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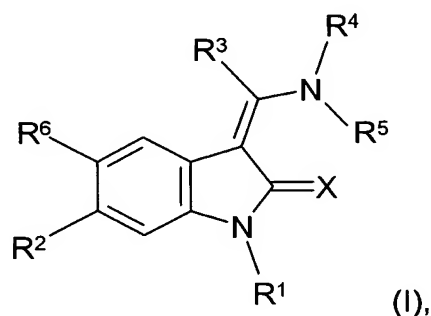
D-55216 Ingelheim

Case 1/1373

Provisional text

Indolinone derivatives substituted in the 6 position, the preparation thereof
and their use as pharmaceutical compositions

The present invention relates to indolinone derivatives substituted in the 6 position of general formula



the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof, particularly the physiologically acceptable salts thereof which have valuable pharmacological properties, pharmaceutical compositions containing these compounds, the use thereof and processes for the preparation thereof.

The above compounds of general formula I have valuable pharmacological properties, particularly an inhibiting effect on various kinases, especially receptor tyrosine kinases such as VEGFR1, VEGFR2, VEGFR3, PDGFR α , PDGFR β , FGFR1, FGFR3, EGFR, HER2, c-Kit, IGF1R and HGFR, Flt-3, as well as on the proliferation of cultivated human cells, particularly endothelial cells, e.g. in angiogenesis, but also on the proliferation of other cells, particularly tumour cells.

The present invention thus relates to the above compounds of general formula I which have valuable pharmacological properties, pharmaceutical compositions containing these pharmacologically active compounds, the use thereof and processes for the preparation thereof.

In the above general formula I

X denotes an oxygen or sulphur atom,

R¹ denotes a hydrogen atom, a C₁₋₄-alkoxy-carbonyl, C₁₋₃-alkyl-carbonyl, aminomethyl, C₁₋₃-alkylaminomethyl, di-(C₁₋₃-alkyl)-aminomethyl or a 5- to 7-membered cycloalkyleneiminomethyl group,

R² denotes a fluorine, chlorine or bromine atom or a cyano group,

R³ denotes a phenyl or naphthyl group or

a phenyl or naphthyl group mono- or disubstituted by a fluorine, chlorine, bromine or iodine atom, by a trifluoromethyl, C₁₋₃-alkyl or C₁₋₃-alkoxy group, while in the case of disubstitution the substituents may be identical or different and the abovementioned unsubstituted as well as the mono- and disubstituted phenyl and naphthyl groups may additionally be substituted

by a fluorine, chlorine, bromine or iodine atom,

by a C₁₋₃-alkyl, C₁₋₄-alkoxy, C₁₋₄-alkoxy-carbonyl, C₁₋₄-alkyloxy-carbonylamino, aminocarbonyl, C₁₋₃-alkylamino-carbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, benzyloxy, carboxy, cyano, trifluoromethyl, nitro, amino, C₄₋₇-cycloalkylamino, C₁₋₃-alkyl-carbonyl-amino, N-(C₁₋₃-alkyl)-N-(C₁₋₃-alkyl-carbonyl)-amino, phenyl-carbonylamino, N-(C₁₋₃-alkyl)-N-(phenyl-carbonyl)-amino, benzyl-carbonylamino, N-(C₁₋₃-alkyl)-N-(benzyl-carbonyl)-amino, hydroxy, C₁₋₃-alkylsulphonylamino, N-(C₁₋₃-alkyl)-C₁₋₃-alkylsulphonylamino, phenylsulphonylamino, N-(C₁₋₃-alkyl)-phenylsulphonylamino, phenyl-

C₁₋₃-alkyl-sulphonylamino, *N*-(C₁₋₃-alkyl)-*N*-(phenyl-C₁₋₃-alkyl-sulphonyl)-amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

by a C₁₋₃-alkyl group which is substituted by a hydroxy, cyano, carboxy, C₁₋₄-alkoxy, C₁₋₄-alkoxy-carbonyl, aminocarbonyl, (C₁₋₃-alkylamino)-carbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, amino, C₁₋₃-alkylamino, [di-(C₁₋₃-alkyl)-amino], *N*-(C₁₋₄-alkoxy-carbonyl)-amino, *N*-(C₁₋₄-alkoxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino, phenylamino, diphenylamino, *N*-phenyl-*N*-(C₁₋₃-alkyl)-amino, benzylamino, dibenzylamino, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino, heteroaryl-amino, *N*-heteroaryl-*N*-(C₁₋₃-alkyl)-amino, C₁₋₄-alkyl-sulphonylamino, *N*-(C₁₋₃-alkyl)-C₁₋₄-alkylsulphonylamino, phenylsulphonylamino, *N*-(C₁₋₃-alkyl)-phenylsulphonylamino, phenyl-C₁₋₃-alkyl-sulphonylamino, *N*-(C₁₋₃-alkyl)-*N*-(phenyl-C₁₋₃-alkyl-sulphonyl)-amino, benzylcarbonylamino, *N*-(C₁₋₃-alkyl)-*N*-(benzylcarbonyl)-amino, phenylcarbonylamino, *N*-(C₁₋₃-alkyl)-*N*-(phenylcarbonyl)-amino, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl, (C₁₋₆-alkyl-carbonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₆-alkyl-carbonyl)-amino, (C₃₋₇-cycloalkyl-carbonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-carbonyl)-amino, (C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino, (C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonyl)-amino, (heteroaryl-carbonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(heteroaryl-carbonyl)-amino, (C₃₋₇-cycloalkyl-sulphonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-sulphonyl)-amino, (C₃₋₇-cycloalkyl-C₁₋₃-alkyl-sulphonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-C₁₋₃-alkyl-sulphonyl)-amino, (C₁₋₄-alkoxy-C₁₋₃-alkyl-sulphonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₄-alkoxy-C₁₋₃-alkyl-sulphonyl)-amino, (heteroaryl-sulphonyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(heteroaryl-sulphonyl)-amino, tetrazolyl or heteroaryl group,

by a carboxy-C₂₋₃-alkenyl, aminocarbonyl-C₂₋₃-alkenyl, (C₁₋₃-alkyl-amino)-carbonyl-C₂₋₃-alkenyl, di-(C₁₋₃-alkyl)-amino-carbonyl-C₂₋₃-alkenyl or C₁₋₄-alkoxy-carbonyl-C₂₋₃-alkenyl group,

by a heteroaryl group or

by a cycloalkyleneimino or cycloalkyleneimino- C_{1-3} -alkyl group with in each case 5 to 7 ring members, wherein in each case a methylene group linked to the imino group is replaced by a carbonyl or sulphonyl group or the two methylene groups linked to the imino group are each replaced by a carbonyl group or a $-CH_2-CH_2-$ group linked to the imino group is replaced by the group $-O-CO-$, while the carbonyl group of the $-O-CO-$ group is linked to the imino group and a phenyl ring may be fused to the 5- to 7-membered cycloalkyleneimino group via two adjacent carbon atoms, or

by a cycloalkyleneimino, cycloalkyleneiminocarbonyl, cycloalkyleneiminosulphonyl, cycloalkyleneimino- C_{1-3} -alkyl, cycloalkyleneiminocarbonyl- C_{1-3} -alkyl or cycloalkyleneiminosulphonyl- C_{1-3} -alkyl group with 4 to 7 ring members in each case, while

in each case the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be substituted by a carboxy, C_{1-4} -alkoxy-carbonyl, aminocarbonyl, C_{1-3} -alkylaminocarbonyl, di- $(C_{1-3}$ -alkyl)-aminocarbonyl, phenyl- C_{1-3} -alkylamino or N- $(C_{1-3}$ -alkyl)-phenyl- C_{1-3} -alkylamino group or

may be replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, $-NH$, $-N(C_{1-3}\text{-alkyl})$, $-N(\text{phenyl})$, $-N(C_{1-3}\text{-alkyl-carbonyl})$ or $-N(\text{benzoyl})$ - group,

while the substituents may be identical or different,

R^4 denotes a benzopyrazolyl group,

a C_{3-7} -cycloalkyl group which may be substituted by a N -[di- $(C_{1-3}$ -alkyl)-amino- C_{1-3} -alkyl-carbonyl]-amino or N -[di- $(C_{1-3}$ -alkyl)-amino- C_{1-3} -alkyl-carbonyl]- N - C_{1-3} -alkyl-amino group,

while the methylene group in the 4 position of a 6- or 7-membered cycloalkyl group may be substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or replaced by a -NH or -N(C₁₋₃-alkyl) group,

or a phenyl, naphthyl or heteroaryl group substituted by the group R₉ which may additionally be mono- or disubstituted by fluorine, chlorine, bromine or iodine atoms, by C₁₋₅-alkyl, trifluoromethyl, hydroxy, C₁₋₄-alkoxy, benzyloxy, carboxy, C₁₋₄-alkoxy-carbonyl, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, acetylamino, C₁₋₃-alkyl-sulphonylamino, aminocarbonyl, C₁₋₃-alkyl-aminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, aminosulphonyl, C₁₋₃-alkyl-aminosulphonyl, di-(C₁₋₃-alkyl)-aminosulphonyl, nitro or cyano groups, while the substituents may be identical or different and wherein

R₉ denotes a hydrogen, fluorine, chlorine, bromine or iodine atom,

a cyano, nitro, amino, C₁₋₅-alkyl, C₃₋₇-cycloalkyl, trifluoromethyl, phenyl, tetrazolyl or heteroaryl group,

a C₁₋₃-alkyl-sulphonyl, amino-C₁₋₃-alkyl-sulphonyl, (C₁₋₃-alkylamino)-C₁₋₃-alkyl-sulphonyl or di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkylsulphonyl group,

a C₁₋₄-alkoxy group, a ω-C₁₋₃-alkoxy-C₂₋₃-alkoxy, phenyl-C₁₋₃-alkoxy, ω-amino-C₂₋₃-alkoxy, ω-(C₁₋₃-alkylamino)-C₂₋₃-alkoxy, ω-[di-(C₁₋₃-alkyl)-amino]-C₂₋₃-alkoxy, ω-(phenyl-C₁₋₃-alkylamino)-C₂₋₃-alkoxy, ω-[N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino]-C₂₋₃-alkoxy, ω-(C₅₋₇-cycloalkyleneimino)-C₂₋₃-alkoxy or C₁₋₃-alkylmercapto group,

a carboxy or C₁₋₄-alkoxy-carbonyl group, aminocarbonyl, C₁₋₄-alkyl-amino-carbonyl, N-(C₁₋₅-alkyl)-C₁₋₃-alkylaminocarbonyl, C₃₋₇-cycloalkyl-amino-carbonyl, N-(C₁₋₅-alkyl)-C₃₋₇-cycloalkylaminocarbonyl, (phenyl-C₁₋₃-alkyl)-amino-carbonyl, N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino-carbonyl group,

a C₁₋₃-alkylaminocarbonyl or *N*-(C₁₋₃-alkyl)-C₁₋₃-alkylaminocarbonyl group, wherein one or two alkyl moieties are substituted independently of one another by a nitro, cyano, carbamoyl, *N*-(C₁₋₃-alkyl)-carbamoyl, di-*N*-(C₁₋₃-alkyl)-carbamoyl, carboxy or C₁₋₄-alkoxy-carbonyl group or in the 2 or 3 position by an amino, (C₁₋₃-alkyl)-amino, di-(C₁₋₃-alkyl)-amino, (C₁₋₄-alkoxy-carbonyl)-amino, *N*-(C₁₋₄-alkoxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino, piperazino, *N*-(C₁₋₃-alkyl)-piperazino, a 4- to 7-membered cycloalkyleneimino group, a hydroxy or methoxy group,

a 4- to 7-membered cycloalkyleneiminocarbonyl group wherein

the cycloalkylene moiety may be fused to a phenyl ring via two adjacent ring atoms or may be bridged to a methylene or ethylene group via two non-adjacent ring atoms or

one or two hydrogen atoms in each case may be replaced by a C₁₋₃-alkyl group and/or

in each case the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneiminocarbonyl group may be substituted by a carboxy, C₁₋₄-alkoxy-carbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-amino, phenyl-C₁₋₃-alkylamino or *N*-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino group, a hydroxy or methoxy group or

replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl or -NH group or by a nitrogen atom, the by a C₁₋₃-alkyl, phenyl, C₁₋₃-alkyl-carbonyl, C₁₋₄-alkoxy-carbonyl, di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, ω-hydroxy-C₂₋₃-alkyl or benzoyl group,

a 4- to 7-membered cycloalkyleneimino group wherein

a methylene group linked to the imino group by a carbonyl or sulphonyl group may be replaced or

the cycloalkylene moiety may be fused to a phenyl ring or

one or two hydrogen atoms in each case may be replaced by a C₁₋₃-alkyl group and/or

in each case the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be substituted by a carboxy, C₁₋₄-alkoxy-carbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, phenyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino group or

replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, -NH, -N(C₁₋₃-alkyl), -N(phenyl), -N(C₁₋₃-alkyl-carbonyl) or -N(benzoyl) group,

a C₁₋₄-alkyl group substituted by the group R₁₀, where

R₁₀ denotes a C₃₋₇-cycloalkyl group,

while the methylene group in the 4 position of a 6- or 7-membered cycloalkyl group may be substituted by an amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or replaced by a -NH or -N(C₁₋₃-alkyl) group or

in a 5- to 7-membered cycloalkyl group a -(CH₂)₂ group may be replaced by a -CO-NH group, a -(CH₂)₃ group may be replaced by a -NH-CO-NH or -CO-NH-CO group or a -(CH₂)₄ group may be replaced by a -NH-CO-NH-CO group, while in each case a hydrogen atom bonded to a nitrogen atom may be replaced by a C₁₋₃-alkyl group,

a phenyl, triazolyl or heteroaryl group,

a hydroxy or C₁₋₄-alkoxy group,

an amino, C₁₋₇-alkylamino, di-(C₁₋₇-alkyl)-amino, phenylamino, *N*-phenyl-*N*-(C₁₋₃-alkyl)-amino, *N*-(phenyl-C₁₋₃-alkyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(phenyl-C₁₋₃-alkyl)-amino or di-(phenyl-C₁₋₃-alkyl)-amino group,

a ω-hydroxy-C₂₋₃-alkyl-amino,
N-(C₁₋₃-alkyl)-(ω-hydroxy-C₂₋₃-alkyl)-amino,
 di-(ω-hydroxy-C₂₋₃-alkyl)-amino or di-(ω-(C₁₋₃-alkoxy)-C₂₋₃-alkyl)-amino-group,

a C₁₋₃-alkyl-carbonylamino-C₂₋₃-alkyl-amino or C₁₋₃-alkyl-carbonyl-amino-C₂₋₃-alkyl-*N*-(C₁₋₃-alkyl)-amino group,

a C₁₋₄-alkyloxy-carbonyl-amino, *N*-(C₁₋₄-alkyloxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino or *N*-{ω-[*N*-(C₁₋₄-alkoxy-carbonyl)-amino]-(C₁₋₄-alkyl)}-*N*-(C₁₋₃-alkyl)-amino group,

a C₁₋₃-alkylsulphonylamino, *N*-(C₁₋₃-alkyl)-C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylsulphonylamino-C₂₋₃-alkyl-amino or C₁₋₃-alkylsulphonyl-amino-C₂₋₃-alkyl-*N*-(C₁₋₃-alkyl)-amino group,

a hydroxycarbonyl-C₁₋₃-alkylamino or *N*-(C₁₋₃-alkyl)-hydroxycarbonyl-C₁₋₃-alkyl-amino group,

an *N*-(ω-amino-C₂₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-(ω-C₁₋₃-alkylamino-C₂₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-[ω-di-(C₁₋₃-alkyl)-amino-C₂₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, *N*-(ω-C₁₋₃-alkoxy-C₂₋₃-alkoxy-C₁₋₃-alkyl)-amino or *N*-(ω-C₁₋₃-alkoxy-C₂₋₃-alkoxy-C₁₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino group,

a guanidino group wherein one or two hydrogen atoms may each be replaced by a C₁₋₃-alkyl group,

a C₄₋₇-cycloalkylamino, C₄₋₇-cycloalkyl-C₁₋₃-alkylamino or C₄₋₇-cyclo-alkenylamino group wherein position 1 of the ring does not participate in the double bond and the abovementioned groups may each additionally be substituted at the aminonitrogen atom by a C₅₋₇-cycloalkyl, C₂₋₄-alkenyl or C₁₋₄-alkyl group,

a 4- to 7-membered cycloalkyleneimino group wherein

the cycloalkylene moiety may be fused with a phenyl group or with an oxazolo, imidazolo, thiazolo, pyridino, pyrazino or pyrimidino group optionally substituted by a fluorine, chlorine, bromine or iodine atom or by a nitro, C₁₋₃-alkyl, C₁₋₃-alkoxy or amino group and/or

one or two hydrogen atoms may each be replaced by a C₁₋₃-alkyl, C₁₋₄-alkoxy-carbonyl, amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group C₅₋₇-cycloalkyl or phenyl group and/or

the methylene group in position 3 of a 5-membered cycloalkyleneimino group may be substituted by a hydroxy, hydroxy-C₁₋₃-alkyl, C₁₋₄-alkoxy or C₁₋₃-alkoxy-C₁₋₃-alkyl group,

in each case the methylene group in position 3 or 4 of a 6- or 7-membered cycloalkyleneimino group may be substituted by a hydroxy, hydroxy-C₁₋₃-alkyl, C₁₋₄-alkoxy, C₁₋₄-alkoxy-C₁₋₃-alkyl, carboxy, C₁₋₄-alkoxy-carbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, *N*-(phenyl-C₁₋₃-alkyl)-amino or *N*-(C₁₋₃-alkyl)-*N*-(phenyl-C₁₋₃-alkyl)-amino group or

may be replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, -NH, -N(C₁₋₃-alkyl-), -N(phenyl), -N(phenyl-C₁₋₃-alkyl-), -N(C₁₋₃-alkyl-carbonyl-), -N(C₁₋₄-hydroxy-carbonyl-),

-N(C₁₋₄-alkoxy-carbonyl-), -N(benzoyl-) or
 -N(phenyl-C₁₋₃-alkyl-carbonyl-) group,

while a methylene group linked to an imino-nitrogen atom of the cycloalkyleneimino group may be replaced by a carbonyl or sulphonyl group or in a 5- to 7-membered monocyclic cycloalkyleneimino group or a cycloalkyleneimino group fused to a phenyl group the two methylene groups linked to the imino-nitrogen atom may each be replaced by a carbonyl group,

or R₉ denotes a C₁₋₄-alkyl group which is substituted by a carboxy, C₁₋₄-alkoxy-carbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl)-aminocarbonyl, N-[amino-C₁₋₃-alkyl]-aminocarbonyl, N-[(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-aminocarbonyl, N-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-aminocarbonyl, N-[amino-C₁₋₃-alkyl]-N-(C₁₋₃-alkyl)-aminocarbonyl, N-[(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-N-(C₁₋₃-alkyl)-aminocarbonyl, N-(C₃₋₇-cycloalkyl)-N-(C₁₋₃-alkyl)-amino or N-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-N-(C₁₋₃-alkyl)-aminocarbonyl group or by a 4- to 7-membered cycloalkyleneiminocarbonyl group,

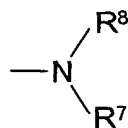
while in the abovementioned cycloalkyleneimino groups one or two hydrogen atoms may each be replaced by a C₁₋₃-alkyl, carboxy, C₁₋₄-alkoxy-carbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl or di-(C₁₋₃-alkyl)-aminocarbonyl group or

one or two hydrogen atoms, which are bonded to a carbon atom not adjacent to the imino group, may be replaced by an amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, phenyl-C₁₋₃-alkylamino or N-(C₁₋₃-alkyl)-phenyl-C₁₋₃-alkylamino group and/or

the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be replaced by one of the groups -S, -SO, -SO₂, -NH, -N(C₁₋₃-alkyl), -N(phenyl), -N(C₁₋₃-alkyl-carbonyl), -N(C₁₋₄-alkoxy-carbonyl), -N(benzoyl) or -O-,

an N-(C₁₋₃-alkyl)-C₁₋₃-alkyl-carbonyl-amino group which is additionally substituted in the alkyl moiety by a carboxy or C₁₋₄-alkoxy-carbonyl group, or

a group of formula



wherein

R⁷ denotes a hydrogen atom, a C₁₋₄-alkyl or C₃₋₇-cycloalkyl group,

a C₁₋₃-alkyl group terminally substituted by a phenyl, heteroaryl, trifluoromethyl, aminocarbonyl, C₁₋₄-alkylamino-carbonyl, di-(C₁₋₄-alkyl)-amino-carbonyl, C₁₋₃-alkyl-carbonyl, C₁₋₃-alkyl-sulphonylamino, N-(C₁₋₃-alkyl)-C₁₋₃-alkyl-sulphonylamino, C₁₋₃-alkyl-aminosulphonyl or di-(C₁₋₃-alkyl)-aminosulphonyl group,

a C₂₋₃-alkyl group terminally substituted by a hydroxy or C₁₋₃-alkoxy group,

a C₁₋₄-alkyl-carbonyl, benzylcarbonyl, heteroarylcarbonyl, heteroaryl-C₁₋₃-alkyl-carbonyl, cycloalkyleneimino-C₁₋₃-alkyl-carbonyl with 5 to 7 ring atoms in the cycloalkyleneimino moiety, C₁₋₃-alkoxy-C₁₋₃-alkyl-carbonyl, amino-C₁₋₃-alkylcarbonyl, (C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-carbonyl, di-(C₁₋₃-alkyl)-amino-carbonyl-C₁₋₃-alkyl, C₁₋₄-alkylsulphonyl, phenylsulphonyl, heteroarylsulphonyl, heteroaryl-C₁₋₃-alkyl-sulphonyl or benzylsulphonyl group or

a phenylcarbonyl group optionally substituted in the phenyl moiety by one or two methoxy groups and

R^8 denotes a C_{1-3} -alkyl, di- $(C_{1-3}$ -alkyl)-amino- C_{1-3} -alkyl-amino-carbonyl or 1- $(C_{1-3}$ -alkyl)-piperidin-4-yl-aminocarbonyl group,

a C_{1-4} -alkyl-carbonyl group terminally substituted by a (ω -alkoxy- C_{2-3} -alkyl)-amino, C_{1-3} -alkyl-carbonyl-amino or *N*-[di- $(C_{1-3}$ -alkyl)-amino- C_{1-3} -alkyl]-*N*- $(C_{1-3}$ -alkyl)-amino group or

a C_{2-4} -alkyl, carbonyl, C_{1-4} -alkyl-carbonyl or carbonyl- C_{1-3} -alkyl group terminally substituted by one of the groups described under R^{10} ,

while R^{10} additionally also denotes a C_{5-7} -cycloalkyloxy group wherein the methylene group may be substituted in the 4 position by a -NH or - $N(C_{1-3}$ -alkyl)- group,

a 5- to 7-membered cycloalkyleneimino-amino group, while the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be substituted by a carboxy, C_{1-3} -alkoxycarbonyl, aminocarbonyl, C_{1-3} -alkylaminocarbonyl, di- $(C_{1-3}$ -alkyl)-aminocarbonyl, phenyl- C_{1-3} -alkylamino or *N*- $(C_{1-3}$ -alkyl)-phenyl- C_{1-3} -alkylamino group or

replaced by an oxygen or sulphur atom, by a sulphinyl, sulphonyl, -NH, - $N(C_{1-3}$ -alkyl), -N(phenyl), - $N(C_{1-3}$ -alkyl-carbonyl) or -N(benzoyl)- group,

or may denote an *N*-(heteroaryl- C_{1-3} -alkyl)-amino group,

R^5 denotes a hydrogen atom or a C_{1-3} -alkyl group and

R^6 denotes a hydrogen atom or a nitro group,

while the unsubstituted, mono- or disubstituted phenyl groups contained in the above definitions, whether singly bonded or fused on, may additionally be substituted by one or two fluorine, chlorine, bromine or iodine atoms or by one or two C_{1-5} -alkyl, C_{1-4} -alkoxy, benzyloxy, carboxy, cyano, C_{1-4} -alkoxy-carbonyl, aminocarbonyl, C_{1-4} -alkylamino-carbonyl, di- $(C_{1-4}$ -alkyl)-amino-carbonyl, aminosulphonyl, C_{1-3} -alkyl-aminosulphonyl, di- $(C_{1-3}$ -alkyl)-aminosulphonyl, trifluoromethyl, nitro, amino, hydroxy, C_{1-3} -alkylsulphonylamino, C_{1-3} -alkylamino or di- $(C_{1-3}$ -alkyl)-amino groups, while the substituents may be identical or different,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while, unless otherwise stated, the expression a heteroaryl group refers to a monocyclic 5- or 6-membered heteroaryl group optionally substituted in the carbon skeleton by a C_{1-3} -alkyl group, wherein

the 6-membered heteroaryl group contains one, two or three nitrogen atoms and

the 5-membered heteroaryl group contains an imino group optionally substituted by a C_{1-3} -alkyl or phenyl- C_{1-3} -alkyl group, an oxygen or sulphur atom or

an imino group optionally substituted by a C_{1-3} -alkyl, amino- C_{1-3} -alkyl, C_{1-3} -alkylamino- C_{1-3} -alkyl, di- $(C_{1-3}$ -alkyl)-amino- C_{1-3} -alkyl or phenyl- C_{1-3} -alkyl group or an oxygen or sulphur atom and additionally contains a nitrogen atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group and two nitrogen atoms,

and moreover a phenyl ring may be fused to the abovementioned monocyclic heterocyclic groups via two adjacent carbon atoms and the bonding takes place via a nitrogen atom or via a carbon atom of the heterocyclic moiety or a fused-on phenyl ring,

and additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved *in vivo*,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof,

with the exception of the compounds

(Z)-3-[1-(4-piperidinomethyl-phenylamino)-1-phenyl-methylidene]-6-chloro-2-indolinone and

(Z)-3-[1-(4-piperidinomethyl-phenylamino)-1-phenyl-methylidene]-6-bromo-2-indolinone.

By a group which may be converted in-vivo into a carboxy group is meant for example a hydroxymethyl group, a carboxy group esterified with an alcohol wherein the alcoholic moiety preferably denotes a C₁₋₆-alkanol, a phenyl-C₁₋₃-alkanol, a C₃₋₉-cycloalkanol, while a C₅₋₈-cycloalkanol may additionally be substituted by one or two C₁₋₃-alkyl groups, a C₅₋₈-cycloalkanol, wherein a methylene group is replaced in the 3 or 4 position by an oxygen atom or by an imino group optionally substituted by a C₁₋₃-alkyl, phenyl-C₁₋₃-alkyl, phenyl-C₁₋₃-alkoxy-carbonyl or C₁₋₆-alkyl-carbonyl group and the cycloalkanol moiety may additionally be substituted by one or two C₁₋₃-alkyl groups, a C₄₋₇-cycloalkenol, a C₃₋₅-alkenol, a phenyl-C₃₋₅-alkenol, a C₃₋₅-alkynol or phenyl-C₃₋₅-alkynol, with the proviso that no bond to the oxygen atom starts from a carbon atom which carries a double or triple bond, a C₃₋₈-cycloalkyl-

C₁₋₃-alkanol, a bicycloalkanol with a total of 8 to 10 carbon atoms which may additionally be substituted in the bicycloalkyl moiety by one or two C₁₋₃-alkyl groups, a 1,3-dihydro-3-oxo-1-isobenzofuranol or an alcohol of formula



wherein

R_a denotes a C₁₋₈-alkyl, C₅₋₇-cycloalkyl, phenyl or phenyl-C₁₋₃-alkyl group,

R_b denotes a hydrogen atom, a C₁₋₃-alkyl, C₅₋₇-cycloalkyl or phenyl group and

R_c denotes a hydrogen atom or a C₁₋₃-alkyl group,

and by a group which can be cleaved *in vivo* from an imino or amino group is meant for example a hydroxy group, an acyl group such as the benzoyl or pyridinoyl group or a C₁₋₁₆-alkyl-carbonyl group such as the formyl, acetyl, propionyl, butanoyl, pentanoyl or hexanoyl group, an allyloxycarbonyl group, a C₁₋₁₆-alkoxy-carbonyl group such as the methoxycarbonyl, ethoxycarbonyl, propoxycarbonyl, isopropoxycarbonyl, butoxycarbonyl, tert.butoxycarbonyl, pentoxycarbonyl, hexyloxycarbonyl, octyloxycarbonyl, nonyloxycarbonyl, decyloxycarbonyl, undecyloxycarbonyl, dodecyloxycarbonyl or hexadecyloxy-carbonyl group, a phenyl-C₁₋₆-alkoxy-carbonyl group such as the benzyloxycarbonyl, phenylethoxycarbonyl or phenylpropoxycarbonyl group, a C₁₋₃-alkylsulphonyl-C₁₋₄-alkoxy-carbonyl, C₁₋₃-alkoxy-C₂₋₄-alkoxy-C₂₋₄-alkoxy-carbonyl or R_aCO-O-(R_bCR_c)-O-CO-group wherein

R_a denotes a C₁₋₈-alkyl, C₅₋₇-cycloalkyl, phenyl or phenyl-C₁₋₃-alkyl group,

R_b denotes a hydrogen atom, a C₁₋₃-alkyl, C₅₋₇-cycloalkyl or phenyl group and

R_c denotes a hydrogen atom, a C_{1-3} -alkyl or $R_aCO-O-(R_bCR_c)-O$ -group wherein R_a to R_c are as hereinbefore defined,

and additionally for an amino group the phthalimido group, while the abovementioned ester groups may also be used as a group which can be converted *in vivo* into a carboxy group.

Preferred prodrug groups for a carboxy group include a C_{1-6} -alkoxy-carbonyl group such as the methoxycarbonyl, ethoxycarbonyl, n-propyloxycarbonyl, isopropyloxycarbonyl, n-butyloxycarbonyl, n-pentyloxycarbonyl, n-hexyloxycarbonyl or cyclohexyloxycarbonyl group or phenyl- C_{1-3} -alkoxy-carbonyl group such as the benzyloxycarbonyl group and

for an imino or amino group a C_{1-9} -alkoxy-carbonyl group such as the methoxycarbonyl, ethoxycarbonyl, n-propyloxycarbonyl, isopropyloxycarbonyl, n-butyloxycarbonyl, n-pentyloxycarbonyl, n-hexyloxycarbonyl, cyclohexyloxycarbonyl, n-heptyloxycarbonyl, n-octyloxycarbonyl or n-nonyloxycarbonyl group, a phenyl- C_{1-3} -alkoxy-carbonyl group such as the benzyloxycarbonyl group, a phenylcarbonyl group optionally substituted by a C_{1-3} -alkyl group such as the benzoyl or 4-ethyl-benzoyl group, a pyridinoyl group such as the nicotinoyl group, a C_{1-3} -alkylsulphonyl-n- C_{2-3} -alkoxy-carbonyl or C_{1-3} -alkoxy- C_{2-3} -alkoxy- C_{1-4} -alkoxy-carbonyl group such as the 2-methylsulphonylethoxycarbonyl or 2-(2-ethoxy)-ethoxycarbonyl group.

Moreover, the saturated alkyl and alkoxy moieties containing more than 2 carbon atoms mentioned in the definitions above as well as the alkanoyl and unsaturated alkyl moieties which contain more than 3 carbon atoms also include the branched isomers thereof such as the isopropyl, tert.butyl, isobutyl group, etc.

One sub-group of compounds of general formula I deserving special mention are those wherein

X denotes an oxygen or sulphur atom,

R^1 denotes a hydrogen atom, a C_{1-4} -alkoxy-carbonyl, C_{1-3} -alkyl-carbonyl, aminomethyl, C_{1-3} -alkylaminomethyl, di- $(C_{1-3}$ -alkyl)-aminomethyl or a 5- to 7-membered cycloalkyleneiminomethyl group,

R^2 denotes a fluorine, chlorine or bromine atom or a cyano group,

R^3 denotes a phenyl or naphthyl group or

a phenyl or naphthyl group mono- or disubstituted by a fluorine, chlorine, bromine or iodine atom, by a trifluoromethyl, C_{1-3} -alkyl or C_{1-3} -alkoxy group, while in the case of disubstitution the substituents may be identical or different and the abovementioned unsubstituted as well as the mono- and disubstituted phenyl and naphthyl groups may additionally be substituted

by a fluorine, chlorine, bromine or iodine atom,

by a C_{1-3} -alkyl, C_{1-4} -alkoxy, benzyloxy, carboxy, cyano, trifluoromethyl, nitro, amino, C_{1-3} -alkyl-carbonyl-amino, C_{1-4} -alkyloxy-carbonylamino, N -(C_{1-3} -alkyl)- N -(C_{1-3} -alkyl-carbonyl)-amino, phenyl-carbonylamino, N -(C_{1-3} -alkyl)- N -(phenyl-carbonyl)-amino, benzyl-carbonylamino, N -(C_{1-3} -alkyl)- N -(benzyl-carbonyl)-amino, hydroxy, C_{1-3} -alkylsulphonylamino, N -(C_{1-3} -alkyl)- N -(C_{1-3} -alkylsulphonyl)-amino, phenylsulphonylamino, N -(C_{1-3} -alkyl)- N -(phenylsulphonyl)-amino, benzylsulphonylamino, N -(C_{1-3} -alkyl)- N -(benzylsulphonyl)-amino, C_{1-3} -alkylamino or di- $(C_{1-3}$ -alkyl)-amino group,

by a hydroxy- C_{1-3} -alkyl, cyano- C_{1-3} -alkyl, carboxy- C_{1-3} -alkyl, C_{1-4} -alkoxy- C_{1-3} -alkyl, amino- C_{1-3} -alkyl, C_{1-3} -alkylamino- C_{1-3} -alkyl, [di- $(C_{1-3}$ -alkyl)-amino]- C_{1-3} -alkyl, benzylamino- C_{1-3} -alkyl, dibenzylamino- C_{1-3} -alkyl, N -benzyl- N -(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl, benzylcarbonylamino- C_{1-3} -alkyl, N -(C_{1-3} -alkyl)- N -(benzylcarbonyl)-amino- C_{1-3} -alkyl, phenylcarbonylamino- C_{1-3} -alkyl, N -(C_{1-3} -alkyl)- N -(phenylcarbonyl)-amino- C_{1-3} -alkyl, phenylamino- C_{1-3} -alkyl, diphenylamino- C_{1-3} -alkyl, N -

phenyl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, heteroaryl-amino-C₁₋₃-alkyl, *N*-heteroaryl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, C₁₋₄-alkyl-sulphonylamino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₄-alkyl-sulphonyl)-amino-C₁₋₃-alkyl, phenyl-sulphonylamino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(phenyl-sulphonyl)-amino-C₁₋₃-alkyl, benzyl-sulphonylamino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(benzyl-sulphonyl)-amino-C₁₋₃-alkyl, C₁₋₄-alkoxy-carbonyl-C₁₋₃-alkyl, *N*-(C₁₋₄-alkoxy-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₄-alkoxy-carbonyl)-amino-C₁₋₃-alkyl, aminocarbonyl-C₁₋₃-alkyl, (C₁₋₃-alkylamino)-carbonyl-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyl, (C₁₋₆-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₆-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₃₋₇-cycloalkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (heteroaryl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(heteroaryl-carbonyl)-amino-C₁₋₃-alkyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl-(C₁₋₃-alkyl), tetrazolyl-C₁₋₃-alkyl or heteroaryl-C₁₋₃-alkyl group,

by a carboxy-C₂₋₃-alkenyl, aminocarbonyl-C₂₋₃-alkenyl, (C₁₋₃-alkyl-amino)-carbonyl-C₂₋₃-alkenyl, di-(C₁₋₃-alkylamino)-carbonyl-C₂₋₃-alkenyl or C₁₋₄-alkoxy-carbonyl-C₂₋₃-alkenyl group or

by a cycloalkyleneimino or cycloalkyleneimino-C₁₋₃-alkyl group each with 5 to 7 ring members, wherein in each case one or two methylene groups adjacent to the nitrogen atom may be replaced by a carbonyl or sulphonyl group or a –CH₂-CH₂- group linked to the imino group may be replaced by the group –O-CO-, while the carbonyl group of the –O-CO- group is linked to the imino group,

while the substituents may be identical or different,

R⁴ denotes a benzopyrazolyl or 1-(C₁₋₃-alkyl)-piperidin-4-yl group,

a cyclohexyl group which is substituted by a *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-carbonyl]-amino or *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-carbonyl]-*N*-C₁₋₃-alkyl-amino group, or

a phenyl, furyl, pyrrolyl, pyridinyl or naphthyl group, each of which is substituted in the carbon skeleton

by a fluorine, chlorine, bromine or iodine atom, by a C₁₋₃-alkyl, C₁₋₄-alkoxy, cyano, nitro, carboxy or trifluoromethyl group,

by a ω-amino-C₂₋₃-alkoxy, ω-[(C₁₋₃-alkyl)-amino]-C₂₋₃-alkoxy, ω-[di-(C₁₋₃-alkyl)-amino]-C₂₋₃-alkoxy, C₁₋₃-alkyl-sulphonyl, (C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-sulphonyl, amino-C₁₋₃-alkyl-sulphonyl, di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-sulphonyl, 4-(C₁₋₃-alkyl)-piperazino or heteroaryl group,

by a C₁₋₃-alkyl group which is terminally substituted by a carboxy, C₁₋₄-alkoxy-carbonyl, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(ω-amino-C₂₋₃-alkyl)-amino, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino, *N*-[ω-(di-(C₁₋₃-alkyl)-amino)-C₂₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino-carbonyl, *N*-(ω-hydroxy-C₂₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, di-(ω-hydroxy-C₂₋₃-alkyl)-amino, *N*-(ω-C₁₋₃-alkoxy-C₂₋₃-alkoxy-C₁₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-(C₁₋₄-alkoxy-carbonyl)-amino, *N*-(C₁₋₄-alkoxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-{ω-[*N*-(C₁₋₄-alkoxy-carbonyl)-amino]-(C₁₋₄-alkyl)}-*N*-(C₁₋₃-alkyl)-amino, heteroaryl, triazolyl or by a 5- to 7-membered cycloalkyleneimino or cycloalkyleneiminocarbonyl group,

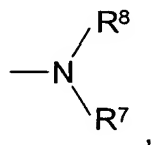
while in the abovementioned cycloalkyleneimino groups one or two hydrogen atoms may each be replaced by a C₁₋₃-alkyl, C₁₋₄-alkoxy-carbonyl, amino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group and/or

the methylene group in the 4 position of a 6- or 7-membered cycloalkyleneimino group may be replaced by one of the groups $-NH$, $-N(C_{1-3}\text{-alkyl})$, $-N(C_{1-4}\text{-alkoxy-carbonyl})$ or $-O-$,

by a carbonyl group which is substituted by a C_{1-3} -alkoxy, N -[amino- C_{1-3} -alkyl]-amino, N -[(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl]-amino, N -[di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl]-amino, N -[amino- C_{1-3} -alkyl]- N -(C_{1-3} -alkyl)-amino, N -[(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl]- N -(C_{1-3} -alkyl)-amino, N -[di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl]- N -(C_{1-3} -alkyl)-amino, N -(C_{3-7} -cycloalkyl)- N -(C_{1-3} -alkyl)-amino or 5- to 7-membered cycloalkyleneimino group,

while the methylene group in the 4 position of a 6- or 7-membered cycloalkylene group may be replaced by a $-NH$, $-N(C_{1-3}\text{-alkyl})$ or $-N(C_{1-4}\text{-alkyloxy-carbonyl})$ -group, or

by a group of formula



wherein

R^7 denotes a hydrogen atom or a C_{1-4} -alkyl, C_{1-4} -alkyl-carbonyl, benzylcarbonyl, heteroarylcarbonyl, cycloalkyleneimino- C_{1-3} -alkyl-carbonyl with 5 to 7 ring atoms in the cycloalkyleneimino moiety, C_{1-3} -alkoxy- C_{1-3} -alkyl-carbonyl, amino- C_{1-3} -alkyl-carbonyl, (C_{1-3} -alkyl)-amino- C_{1-3} -alkyl-carbonyl, di-(C_{1-3} -alkyl)-amino-carbonyl- C_{1-3} -alkyl, C_{1-4} -alkylsulphonyl, phenylsulphonyl, heteroarylsulphonyl or benzylsulphonyl group or a phenylcarbonyl group optionally substituted in the phenyl moiety by one or two methoxy groups and

R⁸ denotes a C₁₋₃-alkyl group, a C₂₋₄-alkyl group terminally substituted by an amino, (C₁₋₃-alkyl)-amino, di-(C₁₋₃-alkyl)-amino or *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino group, an amino-carbonyl-C₁₋₃-alkyl, (C₁₋₃-alkyl)-amino-carbonyl-C₁₋₃-alkyl or di-(C₁₋₃-alkyl)-amino-carbonyl-C₁₋₃-alkyl group,

a di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-amino-carbonyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-aminocarbonyl, 1-(C₁₋₃-alkyl)-piperidin-4-yl-aminocarbonyl, 1-(C₁₋₃-alkyl)-piperidin-4-yl-oxy-carbonyl or (pyridinyl-C₁₋₃-alkyl)-aminocarbonyl group or

a C₁₋₄-alkyl-carbonyl group terminally substituted by a hydroxy, C₁₋₄-alkyloxy, amino, (C₁₋₃-alkyl)-amino, di-(C₁₋₃-alkyl)-amino, (ω-hydroxy-C₂₋₃-alkyl)-amino, di-(ω-hydroxy-C₂₋₃-alkyl)-amino, (ω-alkoxy-C₂₋₃-alkyl)-amino, di-(ω-alkoxy-C₂₋₃-alkyl)-amino, C₁₋₃-alkyl-carbonyl-amino, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, 1-(C₁₋₃-alkyl)-piperidin-4-yl or heteroaryl group or by a 5- to 7-membered cycloalkyleneimino group,

while the cycloalkylene group may be substituted by a C₁₋₃-alkyl group and/or

one or two methylene groups linked to the imino group may be replaced by a carbonyl group and/or

the methylene group in the 4 position of a 6- or 7-membered cycloalkylimino group may be replaced by an -NH, -N(C₁₋₃-alkyl), -N(benzyl), -N(C₁₋₄-alkoxy-carbonyl) or -O- and/or

a phenyl ring may be fused via two adjacent carbon atoms of the cycloalkyleneimino group,

while a 2- or 3-linked pyrrolyl group may additionally be substituted at the nitrogen atom by a C₁₋₃-alkyl group,

R⁵ denotes a hydrogen atom or a C₁₋₃-alkyl group and

R⁶ denotes a hydrogen atom or a nitro group,

while the unsubstituted, mono- or disubstituted phenyl groups contained in the above definitions may additionally be substituted by a fluorine, chlorine, bromine or iodine atom or by a C₁₋₃-alkyl, C₁₋₃-alkoxy, benzyloxy, carboxy, cyano, trifluoromethyl, nitro, amino, hydroxy, C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or by two methyl groups,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while, unless otherwise stated, the expression a heteroaryl group refers to a monocyclic 5- or 6-membered heteroaryl group optionally substituted in the carbon skeleton by a C₁₋₃-alkyl group, wherein

the 6-membered heteroaryl group contains one, two or three nitrogen atoms and

the 5-membered heteroaryl group contains an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group, an oxygen or sulphur atom or

an imino group optionally substituted by a C₁₋₃-alkyl, amino-C₁₋₃-alkyl, [(C₁₋₃-alkyl)-amino]-C₁₋₃-alkyl, [di-(C₁₋₃-alkyl)-amino]-C₁₋₃-alkyl or

phenyl-C₁₋₃-alkyl group or an oxygen or sulphur atom and additionally contains a nitrogen atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group and two nitrogen atoms,

and moreover a phenyl ring may be fused to the abovementioned monocyclic heterocyclic groups via two adjacent carbon atoms and the bonding takes place via a nitrogen atom or via a carbon atom of the heterocyclic moiety or a fused-on phenyl ring,

and additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved *in vivo*,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof,

with the exception of the compounds

(Z)-3-[1-(4-piperidinomethyl-phenylamino)-1-phenyl-methylidene]-6-chloro-2-indolinone and

(Z)-3-[1-(4-piperidinomethyl-phenylamino)-1-phenyl-methylidene]-6-bromo-2-indolinone.

Preferred compounds of the above general formula I are those wherein

X denotes an oxygen or sulphur atom,

R¹ denotes a hydrogen atom, a C₁₋₄-alkoxy-carbonyl, C₁₋₃-alkyl-carbonyl, aminomethyl, C₁₋₃-alkylaminomethyl, di-(C₁₋₃-alkyl)-aminomethyl or a 5- to 7-membered cycloalkyleneiminomethyl group,

R² denotes a fluorine, chlorine or bromine atom or a cyano group,

R³ denotes a phenyl or naphthyl group or

a phenyl or naphthyl group mono- or disubstituted by a fluorine, chlorine, bromine or iodine atom, by a trifluoromethyl, C₁₋₃-alkyl or C₁₋₃-alkoxy group, while in the case of disubstitution the substituents may be identical or different and the abovementioned unsubstituted as well as the mono- and disubstituted phenyl and naphthyl groups may additionally be substituted

by a fluorine, chlorine, bromine or iodine atom,

by a C₁₋₃-alkyl, C₁₋₄-alkoxy, benzyloxy, carboxy, cyano, trifluoromethyl, nitro, amino, C₁₋₃-alkyl-carbonyl-amino, C₁₋₄-alkyloxy-carbonylamino, hydroxy, C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

by a hydroxy-C₁₋₃-alkyl, cyano-C₁₋₃-alkyl, carboxy-C₁₋₃-alkyl, C₁₋₄-alkoxy-C₁₋₃-alkyl, amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl, [di-(C₁₋₃-alkyl)-amino]-C₁₋₃-alkyl, benzylamino-C₁₋₃-alkyl, dibenzylamino-C₁₋₃-alkyl, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, benzylcarbonylamino-C₁₋₃-alkyl, phenylcarbonylamino-C₁₋₃-alkyl, phenylamino-C₁₋₃-alkyl, diphenylamino-C₁₋₃-alkyl, *N*-phenyl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, heteroaryl-amino-C₁₋₃-alkyl, *N*-heteroaryl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, C₁₋₄-alkyl-sulphonylamino-C₁₋₃-alkyl, phenyl-sulphonylamino-C₁₋₃-alkyl, benzyl-sulphonylamino-C₁₋₃-alkyl, C₁₋₄-alkoxy-carbonyl-C₁₋₃-alkyl, *N*-(C₁₋₄-alkoxy-carbonyl)-amino-C₁₋₃-alkyl, aminocarbonyl-C₁₋₃-alkyl, (C₁₋₃-alkylamino)-carbonyl-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyl, (C₁₋₆-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₆-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₃₋₇-cycloalkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (heteroaryl-carbonyl)-amino-C₁₋₃-alkyl, *N*-

(C₁₋₃-alkyl)-*N*-(heteroaryl-carbonyl)-amino-C₁₋₃-alkyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl-(C₁₋₃-alkyl), tetrazolyl-C₁₋₃-alkyl or imidazolyl-C₁₋₃-alkyl group,

by a carboxy-C₂₋₃-alkenyl, aminocarbonyl-C₂₋₃-alkenyl or C₁₋₄-alkoxy-carbonyl-C₂₋₃-alkenyl group or

by a 5- to 7-membered cycloalkyleneimino group wherein one or two methylene groups adjacent to the nitrogen atom may be replaced by a carbonyl group or a –CH₂-CH₂- group linked to the imino group may be replaced by the group –O-CO-, while the carbonyl group of the –O-CO-group is linked to the imino group,

while the substituents may be identical or different,

R⁴ denotes a benzopyrazolyl or 1-(C₁₋₃-alkyl)-piperidin-4-yl group,

a cyclohexyl group which is substituted by a *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-carbonyl]-amino or *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-carbonyl]-*N*-(C₁₋₃-alkyl)-amino group, or

a phenyl, furyl, pyrrolyl, pyridinyl or naphthyl group each of which may be substituted in the carbon skeleton

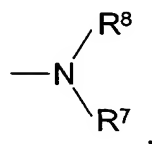
by a fluorine, chlorine, bromine or iodine atom, by a C₁₋₃-alkyl, C₁₋₃-alkoxy, cyano, nitro, carboxy or trifluoromethyl group,

by a ω-[di-(C₁₋₃-alkyl)-amino]-C₂₋₃-alkoxy, C₁₋₃-alkyl-sulphonyl, di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-sulphonyl, 4-(C₁₋₃-alkyl)-piperazino, imidazolyl, C₁₋₃-alkyl-imidazolyl or [di-(C₁₋₃-alkyl)-amino]-C₁₋₃-alkyl-imidazolyl group,

by a C₁₋₃-alkyl group which is terminally substituted by a carboxy, C₁₋₄-alkoxy-carbonyl, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(ω-amino-C₂₋₃-alkyl)-amino, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino, *N*-[ω-(di-(C₁₋₃-alkyl)-amino)-C₂₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino-carbonyl, *N*-(ω-hydroxy-C₂₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, di-(ω-hydroxy-C₂₋₃-alkyl)-amino, *N*-(ω-C₁₋₃-alkoxy-C₂₋₃-alkoxy-C₁₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-(C₁₋₄-alkoxy-carbonyl)-amino, *N*-(C₁₋₄-alkoxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-{ω-[*N*-(C₁₋₄-alkoxy-carbonyl)-amino]-(C₁₋₄-alkyl)}-*N*-(C₁₋₃-alkyl)-amino, pyridinyl, triazolyl, pyrrolidino, piperidino, di-(C₁₋₃-alkyl)-piperidino, [di-(C₁₋₃-alkyl)-amino]-piperidino, piperazino, morpholino, (C₁₋₃-alkyl)-piperazino, (C₁₋₃-alkyl)-piperazin-1-yl-carbonyl or 4-(C₁₋₄-alkoxy-carbonyl)-piperazino group,

by a carbonyl group which is substituted by a C₁₋₄-alkoxy, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, *N*-(C₃₋₇-cycloalkyl)-*N*-(C₁₋₃-alkyl)-amino, piperidino, piperazino, 4-(C₁₋₄-alkoxy-carbonyl)-piperazino or 4-(C₁₋₃-alkyl)-piperazino group, or

by a group of formula



wherein

R⁷ denotes a hydrogen atom or a C₁₋₃-alkyl, C₁₋₃-alkyl-carbonyl, benzylcarbonyl, pyridinylcarbonyl, furanylcarbonyl, pyrrolidino-C₁₋₃-alkyl-carbonyl, C₁₋₃-alkoxy-C₁₋₃-alkyl-carbonyl, di-(C₁₋₃-alkyl)-amino-carbonyl-C₁₋₃-alkyl, C₁₋₄-alkylsulphonyl or benzylsulphonyl

group or a phenylcarbonyl group optionally substituted in the phenyl moiety by one or two methoxy groups and

R^8 denotes a C_{1-3} -alkyl, di- $(C_{1-3}$ -alkyl)-amino- C_{1-4} -alkyl-amino-carbonyl or di- $(C_{1-3}$ -alkyl)-amino-carbonyl- C_{1-3} -alkyl group,

a C_{2-3} -alkyl group terminally substituted by a di- $(C_{1-3}$ -alkyl)-amino or *N*-benzyl-*N*- $(C_{1-3}$ -alkyl)-amino group,

a 4- $(C_{1-3}$ -alkyl)-piperazin-1-yl-carbonyl, 4- $(C_{1-3}$ -alkyl)-piperazin-1-yl-aminocarbonyl, 1- $(C_{1-3}$ -alkyl)-piperidin-4-yl-aminocarbonyl, 1- $(C_{1-3}$ -alkyl)-piperidin-4-yl-oxy-carbonyl or (pyridinyl- C_{1-3} -alkyl)-aminocarbonyl group or

a C_{1-4} -alkyl-carbonyl group terminally substituted by a hydroxy, amino, di- $(C_{1-3}$ -alkyl)-amino, di-(ω -hydroxy- C_{2-3} -alkyl)-amino, C_{1-3} -alkyl-carbonyl-amino, *N*-benzyl-*N*- $(C_{1-3}$ -alkyl)-amino, *N*-[di- $(C_{1-3}$ -alkyl)-amino- C_{1-3} -alkyl]-*N*- $(C_{1-3}$ -alkyl)-amino, imidazolyl, piperazino, 4- $(C_{1-3}$ -alkyl)-piperazin-1-yl, 4-benzyl-piperazin-1-yl, 4- $(C_{1-4}$ -alkoxy-carbonyl)-piperazin-1-yl, 4- $(C_{1-3}$ -alkyl)-homo-piperazin-1-yl, morpholino, pyrrolidino, piperidino, 1- $(C_{1-3}$ -alkyl)-piperidin-4-yl, 4- $(C_{1-3}$ -alkyl)-piperidin-1-yl or phthalimido group,

while a 2- or 3-linked pyrrolyl group may additionally be substituted at the nitrogen atom by a C_{1-3} -alkyl group,

R^5 denotes a hydrogen atom or a C_{1-3} -alkyl group and

R^6 denotes a hydrogen atom or a nitro group,

while the unsubstituted, mono- or disubstituted phenyl groups contained in the above definitions may additionally be substituted by a fluorine, chlorine, bromine or iodine atom or by a C_{1-3} -alkyl, C_{1-3} -alkoxy, benzyloxy, carboxy,

cyano, trifluoromethyl, nitro, amino, hydroxy, C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or by two methyl groups,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while, unless otherwise stated, the expression a heteroaryl group refers to a monocyclic 5- or 6-membered heteroaryl group optionally substituted in the carbon skeleton by a C₁₋₃-alkyl group, wherein

the 6-membered heteroaryl group contains one, two or three nitrogen atoms and

the 5-membered heteroaryl group contains an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group, an oxygen or sulphur atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group or an oxygen or sulphur atom and additionally a nitrogen atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group and two nitrogen atoms,

and moreover a phenyl ring may be fused to the abovementioned monocyclic heterocyclic groups via two adjacent carbon atoms and the bonding takes place via a nitrogen atom or via a carbon atom of the heterocyclic moiety or a fused-on phenyl ring,

and additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved *in vivo*,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof,

with the exception of the compounds

(Z)-3-[1-(4-piperidinomethyl-phenylamino)-1-phenyl-methylidene]-6-chloro-2-indolinone and

(Z)-3-[1-(4-piperidinomethyl-phenylamino)-1-phenyl-methylidene]-6-bromo-2-indolinone.

A first sub-group of preferred compounds of general formula I deserving special mention comprises those wherein

R^1 , R^2 , R^3 , R^5 , R^6 and X are as hereinbefore defined and

R^4 denotes a benzopyrazolyl or 1-(C_{1-3} -alkyl)-piperidin-4-yl group,

a cyclohexyl group which is substituted by an *N*-[di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl-carbonyl]-amino or *N*-[di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl-carbonyl]-*N*- C_{1-3} -alkyl-amino group, or

a phenyl, furyl, pyrrolyl, pyridinyl or naphthyl group, each of which may be substituted in the carbon skeleton

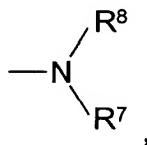
by a fluorine, chlorine, bromine or iodine atom, by a C_{1-3} -alkyl, C_{1-3} -alkoxy, cyano, nitro, carboxy or trifluoromethyl group,

by a ω -[di-(C_{1-3} -alkyl)-amino]- C_{2-3} -alkoxy, C_{1-3} -alkyl-sulphonyl, di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl-sulphonyl, 4-(C_{1-3} -alkyl)-piperazino, imidazolyl, C_{1-3} -alkyl-imidazolyl or [di-(C_{1-3} -alkyl)-amino]- C_{1-3} -alkyl-imidazolyl group,

by a carbonyl group which is substituted by a C_{1-4} -alkoxy, *N*-[di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl]-amino, *N*-[di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl]-*N*-

C₁₋₃-alkyl-amino, *N*-(C₃₋₇-cycloalkyl)-*N*-(C₁₋₃-alkyl)-amino, piperidino, piperazino, 4-(C₁₋₄-alkyloxy-carbonyl)-piperazino or 4-(C₁₋₃-alkyl)-piperazino group, or

by a group of formula



wherein

R⁷ denotes a hydrogen atom or a C₁₋₃-alkyl, C₁₋₃-alkyl-carbonyl, benzylcarbonyl, pyridinylcarbonyl, furanylcabonyl, pyrrolidino-C₁₋₃-alkyl-carbonyl, C₁₋₃-alkoxy-C₁₋₃-alkyl-carbonyl, di-(C₁₋₃-alkyl)-amino-carbonyl-C₁₋₃-alkyl, C₁₋₄-alkylsulphonyl or benzylsulphonyl group or a phenylcarbonyl group optionally substituted in the phenyl moiety by one or two methoxy groups and

R⁸ denotes a C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-amino-carbonyl or di-(C₁₋₃-alkyl)-amino-carbonyl-C₁₋₃-alkyl group,

a C₂₋₃-alkyl group terminally substituted by a di-(C₁₋₃-alkyl)-amino or *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino group,

a 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-aminocarbonyl, 1-(C₁₋₃-alkyl)-piperidin-4-yl-aminocarbonyl, 1-(C₁₋₃-alkyl)-piperidin-4-yl-oxy-carbonyl or (pyridinyl-C₁₋₃-alkyl)-aminocarbonyl group or

a C₁₋₄-alkyl-carbonyl group terminally substituted by a hydroxy, amino, di-(C₁₋₃-alkyl)-amino, di-(ω-hydroxy-C₂₋₃-alkyl)-amino, C₁₋₃-alkyl-carbonyl-amino, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino, *N*-[di-

(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, imidazolyl, piperazino, 4-(C₁₋₃-alkyl)-piperazin-1-yl, 4-benzyl-piperazin-1-yl, 4-(C₁₋₄-alkoxy-carbonyl)-piperazin-1-yl, 4-(C₁₋₃-alkyl)-homo-piperazin-1-yl, morpholino, pyrrolidino, piperidino, 1-(C₁₋₃-alkyl)-piperidin-4-yl, 4-(C₁₋₃-alkyl)-piperidin-1-yl or phthalimido group,

while a 2- or 3-linked pyrrolyl group may additionally be substituted at the nitrogen atom by a C₁₋₃-alkyl group,

while the unsubstituted, mono- or disubstituted phenyl groups contained in the above definitions may additionally be substituted by a fluorine, chlorine, bromine or iodine atom or by a C₁₋₃-alkyl, C₁₋₃-alkoxy, benzyloxy, carboxy, cyano, trifluoromethyl, nitro, amino, hydroxy, C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or by two methyl groups,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while, unless otherwise stated, the expression a heteroaryl group refers to a monocyclic 5- or 6-membered heteroaryl group optionally substituted in the carbon skeleton by a C₁₋₃-alkyl group, wherein

the 6-membered heteroaryl group contains one, two or three nitrogen atoms and

the 5-membered heteroaryl group contains an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group, an oxygen or sulphur atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group or an oxygen or sulphur atom and additionally a nitrogen atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group and two nitrogen atoms,

and moreover a phenyl ring may be fused to the abovementioned monocyclic heterocyclic groups via two adjacent carbon atoms and the bonding takes place via a nitrogen atom or via a carbon atom of the heterocyclic moiety or a fused-on phenyl ring,

and additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved *in vivo*,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

A second sub-group of preferred compounds of general formula I deserving special mention are those wherein

R¹, R², R⁴, R⁵, R⁶ and X are as hereinbefore defined and

R³ denotes a phenyl or naphthyl group or

a phenyl or naphthyl group mono- or disubstituted by a fluorine, chlorine, bromine or iodine atom, by a trifluoromethyl, C₁₋₃-alkyl or C₁₋₃-alkoxy group, while in the case of disubstitution the substituents may be identical or different and the abovementioned unsubstituted as well as the mono- and disubstituted phenyl and naphthyl groups may additionally be substituted

by a fluorine, chlorine, bromine or iodine atom,

by a C₁₋₃-alkyl, C₁₋₄-alkoxy, benzyloxy, carboxy, cyano, trifluoromethyl, nitro, amino, C₁₋₃-alkyl-carbonyl-amino, C₁₋₄-alkyloxy-carbonylamino, hydroxy, C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group,

by a hydroxy-C₁₋₃-alkyl, cyano-C₁₋₃-alkyl, carboxy-C₁₋₃-alkyl, C₁₋₄-alkoxy-C₁₋₃-alkyl, amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl, [di-(C₁₋₃-alkyl)-amino]-C₁₋₃-alkyl, benzylamino-C₁₋₃-alkyl, dibenzylamino-C₁₋₃-alkyl, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, benzylcarbonylamino-C₁₋₃-alkyl, phenylcarbonylamino-C₁₋₃-alkyl, phenylamino-C₁₋₃-alkyl, diphenylamino-C₁₋₃-alkyl, *N*-phenyl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, heteroaryl-amino-C₁₋₃-alkyl, *N*-heteroaryl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, C₁₋₄-alkyl-sulphonylamino-C₁₋₃-alkyl, phenyl-sulphonylamino-C₁₋₃-alkyl, benzyl-sulphonylamino-C₁₋₃-alkyl, C₁₋₄-alkoxy-carbonyl-C₁₋₃-alkyl, *N*-(C₁₋₄-alkoxy-carbonyl)-amino-C₁₋₃-alkyl, aminocarbonyl-C₁₋₃-alkyl, (C₁₋₃-alkylamino)-carbonyl-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyl, (C₁₋₆-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₆-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₃₋₇-cycloalkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonyl)-amino-C₁₋₃-alkyl, (heteroaryl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkyl)-*N*-(heteroaryl-carbonyl)-amino-C₁₋₃-alkyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl-(C₁₋₃-alkyl), tetrazolyl-C₁₋₃-alkyl or imidazolyl-C₁₋₃-alkyl group,

by a carboxy-C₂₋₃-alkenyl, aminocarbonyl-C₂₋₃-alkenyl or C₁₋₄-alkoxy-carbonyl-C₂₋₃-alkenyl group or

by a 5- to 7-membered cycloalkyleneimino group wherein one or two methylene groups adjacent to the nitrogen atom may be replaced by a carbonyl group or a -CH₂-CH₂- group linked to the imino group may be replaced by the group -O-CO-, while the carbonyl group of the -O-CO- group is linked to the imino group,

while the substituents may be identical or different,

and the unsubstituted, mono- or disubstituted phenyl groups contained in the above definitions may additionally be substituted by a fluorine, chlorine, bromine or iodine atom or by a C₁₋₃-alkyl, C₁₋₃-alkoxy, benzyloxy, carboxy, cyano, trifluoromethyl, nitro, amino, hydroxy, C₁₋₃-alkylsulphonylamino, C₁₋₃-alkylamino or di-(C₁₋₃-alkyl)-amino group or by two methyl groups,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while, unless otherwise stated, the expression a heteroaryl group refers to a monocyclic 5- or 6-membered heteroaryl group optionally substituted in the carbon skeleton by a C₁₋₃-alkyl group, wherein

the 6-membered heteroaryl group contains one, two or three nitrogen atoms and

the 5-membered heteroaryl group contains an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group, an oxygen or sulphur atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group or an oxygen or sulphur atom and additionally a nitrogen atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group and two nitrogen atoms,

and moreover a phenyl ring may be fused to the abovementioned monocyclic heterocyclic groups via two adjacent carbon atoms and the bonding takes place via a nitrogen atom or via a carbon atom of the heterocyclic moiety or a fused-on phenyl ring,

and additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved *in vivo*,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

Particularly preferred compounds of the above general formula I are those wherein

X denotes an oxygen or sulphur atom,

R¹ denotes a hydrogen atom, a C₁₋₄-alkoxy-carbonyl, C₁₋₃-alkyl-carbonyl, aminomethyl, C₁₋₃-alkylaminomethyl, di-(C₁₋₃-alkyl)-aminomethyl or a 5- to 7-membered cycloalkyleneiminomethyl group,

R² denotes a fluorine, chlorine or bromine atom or a cyano group,

R³ denotes a phenyl or naphthyl group or

a phenyl or naphthyl group mono- or disubstituted by a fluorine, chlorine, bromine or iodine atom, by a trifluoromethyl, C₁₋₃-alkyl or C₁₋₃-alkoxy group, while in the case of disubstitution the substituents may be identical or different and the abovementioned unsubstituted as well as the mono- and disubstituted phenyl and naphthyl groups may additionally be substituted

by a C₁₋₃-alkyl-carbonyl-amino, C₁₋₄-alkyloxy-carbonylamino, benzyloxy or hydroxy group,

by a hydroxy-C₁₋₃-alkyl, carboxy-C₁₋₃-alkyl, C₁₋₄-alkoxy-C₁₋₃-alkyl, cyano-C₁₋₃-alkyl, benzylamino-C₁₋₃-alkyl, dibenzylamino-C₁₋₃-alkyl, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, C₁₋₄-alkoxy-C₁₋₃-alkyl-carbonylamino-C₁₋₃-alkyl, C₃₋₇-cycloalkyl-carbonylamino-C₁₋₃-alkyl, C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonylamino-C₁₋₃-alkyl, benzylcarbonylamino-C₁₋₃-alkyl, phenylcarbonylamino-C₁₋₃-alkyl, heteroaryl-carbonylamino-C₁₋₃-alkyl, phenylamino-C₁₋₃-alkyl, diphenylamino-C₁₋₃-alkyl, *N*-phenyl-*N*-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, heteroaryl-amino-C₁₋₃-alkyl, *N*-heteroaryl-

N-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl, C₁₋₄-alkyl-sulphonylamino-C₁₋₃-alkyl, phenyl-sulphonylamino-C₁₋₃-alkyl, benzyl-sulphonylamino-C₁₋₃-alkyl, C₁₋₄-alkoxy-carbonyl-C₁₋₃-alkyl, *N*-(C₁₋₄-alkoxy-carbonyl)-amino-C₁₋₃-alkyl, aminocarbonyl-C₁₋₃-alkyl, (C₁₋₃-alkylamino)-carbonyl-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl-(C₁₋₃-alkyl) or tetrazolyl-C₁₋₃-alkyl group,

by an aminocarbonyl-C₂₋₃-alkenyl or C₁₋₄-alkoxy-carbonyl-C₂₋₃-alkenyl group or

by a 5- to 7-membered cycloalkyleneimino groups, wherein one or two methylene groups adjacent to the nitrogen atom may be replaced by a carbonyl group or a -CH₂-CH₂- group linked to the imino group may be replaced by the group -O-CO-, while the carbonyl group of the -O-CO- group is linked to the imino group,

while the substituents may be identical or different,

R⁴ denotes a benzopyrazolyl or 1-(C₁₋₃-alkyl)-piperidin-4-yl group,

a cyclohexyl group which is substituted by a *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-carbonyl]-amino or *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-carbonyl]-*N*-(C₁₋₃-alkyl)-amino group, or

a phenyl, pyridinyl or naphthyl group or a pyrrolyl group optionally substituted at the nitrogen by a C₁₋₃-alkyl group, each of which may be substituted

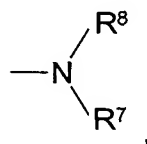
by a fluorine, chlorine, bromine or iodine atom, by a C₁₋₃-alkyl, C₁₋₃-alkoxy, cyano, nitro, carboxy or trifluoromethyl group,

by a ω-[di-(C₁₋₃-alkyl)-amino]-C₂₋₃-alkoxy, C₁₋₃-alkyl-sulphonyl, di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl-sulphonyl, 4-(C₁₋₃-alkyl)-piperazino, imidazolyl, C₁₋₃-alkylimidazolyl or [di-(C₁₋₃-alkyl)-amino]-C₁₋₃-alkyl-imidazolyl group,

by a C₁₋₃-alkyl group which is terminally substituted by a carboxy, C₁₋₃-alkoxy-carbonyl, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, *N*-(C₁₋₃-alkyl)-*N*-(ω-amino-C₂₋₃-alkyl)-amino, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino, *N*-[ω-(di-(C₁₋₃-alkyl)-amino)-C₂₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino-carbonyl, *N*-(ω-hydroxy-C₂₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, di-(ω-hydroxy-C₂₋₃-alkyl)-amino, *N*-(ω-C₁₋₃-alkoxy-C₂₋₃-alkoxy-C₁₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-(C₁₋₄-alkoxy-carbonyl)-amino, *N*-(C₁₋₄-alkoxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-{ω-[*N*-(C₁₋₄-alkoxy-carbonyl)-amino]-(C₁₋₃-alkyl)}-*N*-(C₁₋₃-alkyl)-amino, pyridinyl, triazolyl, pyrrolidino, piperidino, di-(C₁₋₃-alkyl)-piperidino, [di-(C₁₋₃-alkyl)-amino]-piperidino, piperazino, morpholino, (C₁₋₃-alkyl)-piperazino, 4-(C₁₋₃-alkyl)-piperazino-carbonyl or 4-(C₁₋₄-alkoxy-carbonyl)-piperazino group,

by a carbonyl group which is substituted by a C₁₋₃-alkoxy, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, *N*-(C₃₋₇-cycloalkyl)-*N*-(C₁₋₃-alkyl)-amino, piperidino, piperazino, 4-(C₁₋₄-alkyloxy-carbonyl)-piperazino or (C₁₋₃-alkyl)-piperazino group, or

by a group of formula



wherein

R⁷ denotes a hydrogen atom, a C₁₋₃-alkyl, C₁₋₃-alkyl-carbonyl, benzylcarbonyl, pyridinylcarbonyl, furanylcabonyl, pyrrolidino-C₁₋₃-alkyl-carbonyl, C₁₋₃-alkoxy-C₁₋₃-alkyl-carbonyl, di-(C₁₋₃-alkyl)-amino-carbonyl-C₁₋₃-alkyl, C₁₋₄-alkylsulphonyl or benzylsulphonyl group or a phenylcarbonyl group optionally substituted in the phenyl moiety by one or two methoxy groups and

R^8 denotes a C_{1-3} -alkyl, ω -[di-(C_{1-3} -alkyl)-amino]- C_{2-3} -alkyl, di-(C_{1-3} -alkyl)-amino-carbonyl- C_{1-3} -alkyl, di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl-amino-carbonyl or ω -[*N*-benzyl-*N*-(C_{1-3} -alkyl)-amino]- C_{2-3} -alkyl group or

a 4-(C_{1-3} -alkyl)-piperazin-1-yl-carbonyl, 4-(C_{1-3} -alkyl)-piperazin-1-yl-aminocarbonyl, 1-(C_{1-3} -alkyl)-piperidin-4-yl-aminocarbonyl, 1-(C_{1-3} -alkyl)-piperidin-4-yl-oxy-carbonyl or (pyridinyl- C_{1-3} -alkyl)-aminocarbonyl group or

a C_{1-3} -alkyl-carbonyl group terminally substituted by a hydroxy, amino, di-(C_{1-3} -alkyl)-amino, di-(ω -hydroxy- C_{2-3} -alkyl)-amino, C_{1-3} -alkyl-carbonyl-amino, *N*-benzyl-*N*-(C_{1-3} -alkyl)-amino, *N*-[di-(C_{1-3} -alkyl)-amino- C_{1-3} -alkyl]-*N*-(C_{1-3} -alkyl)-amino, imidazolyl, piperazino, 4-(C_{1-3} -alkyl)-piperazin-1-yl, 4-benzyl-piperazin-1-yl, 4-(C_{1-4} -alkoxy-carbonyl)-piperazin-1-yl, 4-(C_{1-3} -alkyl)-homopiperazin-1-yl, morpholino, pyrrolidino, piperidino-, 1-(C_{1-3} -alkyl)-piperidin-4-yl, 4-(C_{1-3} -alkyl)-piperidin-1-yl or phthalimido group,

R^5 denotes a hydrogen atom or a C_{1-3} -alkyl group and

R^6 denotes a hydrogen atom or a nitro group,

while the unsubstituted, mono- or disubstituted phenyl groups contained in the above definitions may additionally be substituted by a cyano or a methoxy group or by two methyl groups, and

the abovementioned alkyl groups include straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while, unless otherwise stated, the expression a heteroaryl group refers to a monocyclic 5- or 6-membered heteroaryl group optionally substituted in the carbon skeleton by a C₁₋₃-alkyl group, wherein

the 6-membered heteroaryl group contains one, two or three nitrogen atoms and

the 5-membered heteroaryl group contains an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group, an oxygen or sulphur atom or

an imino group optionally substituted by a C₁₋₃-alkyl, amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group or an oxygen or sulphur atom and additionally a nitrogen atom or

an imino group optionally substituted by a C₁₋₃-alkyl or phenyl-C₁₋₃-alkyl group and two nitrogen atoms,

and moreover a phenyl ring may be fused to the abovementioned monocyclic heterocyclic groups via two adjacent carbon atoms and the bonding takes place via a nitrogen atom or via a carbon atom of the heterocyclic moiety or a fused-on phenyl ring,

and additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved *in vivo*,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

Most particularly preferred compounds of the above general formula I are those wherein

X denotes an oxygen atom,

R^1 denotes a hydrogen atom,

R^2 denotes a fluorine, chlorine or bromine atom or a cyano group,

R^3 denotes a phenyl group or a phenyl group monosubstituted by a fluorine, chlorine, bromine or iodine atom or by a C_{1-3} -alkoxy group, while the abovementioned unsubstituted and monosubstituted phenyl groups may additionally be substituted in the 3 or 4-position

by a fluorine, chlorine or bromine atom,

by a C_{1-3} -alkoxy or C_{1-2} -alkyl-carbonyl-amino group,

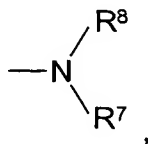
by a carboxy- C_{1-3} -alkyl, aminocarbonyl- C_{1-3} -alkyl, (C_{1-2} -alkylamino)-carbonyl- C_{1-3} -alkyl, di-(C_{1-2} -alkyl)-aminocarbonyl- C_{1-3} -alkyl, (C_{1-2} -alkyl-carbonyl)-amino- C_{1-3} -alkyl or (phenyl-carbonyl)-amino- C_{1-3} -alkyl group,

while the substituents may be identical or different,

R^4 denotes a phenyl group which is substituted

by a C_{1-3} -alkyl group terminally substituted by a di-(C_{1-2} -alkyl)-amino group or

by a group of formula



wherein

R^7 denotes a C_{1-2} -alkyl, C_{1-2} -alkyl-carbonyl, di-(C_{1-2} -alkyl)-aminocarbonyl- C_{1-3} -alkyl or C_{1-3} -alkylsulphonyl group and

R^8 denotes a C_{1-3} -alkyl or ω -[di-(C_{1-2} -alkyl)-amino]- C_{2-3} -alkyl group or

a C_{1-3} -alkyl-carbonyl group terminally substituted by a di-(C_{1-2} -alkyl)-amino, piperazino or 4-(C_{1-3} -alkyl)-piperazin-1-yl group,

R^5 denotes a hydrogen atom and

R^6 denotes a hydrogen atom,

while the abovementioned alkyl groups include straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved *in vivo*,

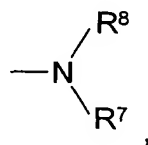
the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

A first sub-group of most particularly preferred compounds of general formula I deserving special mention comprises those wherein

X, R^1 , R^2 , R^3 , R^5 and R^6 are as hereinbefore defined and

R^4 denotes a phenyl group which is substituted

by a group of formula



wherein

R^7 denotes a C_{1-2} -alkyl, C_{1-2} -alkyl-carbonyl, di- $(C_{1-2}$ -alkyl)-amino-carbonyl- C_{1-3} -alkyl or C_{1-3} -alkylsulphonyl group and

R^8 denotes a C_{1-3} -alkyl or ω -[di- $(C_{1-2}$ -alkyl)-amino]- C_{2-3} -alkyl group or

a C_{1-3} -alkyl-carbonyl group terminally substituted by a di- $(C_{1-2}$ -alkyl)-amino, piperazino or 4- $(C_{1-3}$ -alkyl)-piperazin-1-yl group,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved in vivo,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

A second sub-group of particularly preferred compounds of general formula I deserving special mention comprises those wherein

X, R^1 , R^2 , R^4 , R^5 and R^6 are as hereinbefore defined and

R^3 denotes a phenyl group or a phenyl group monosubstituted by a fluorine, chlorine or bromine atom or by a C_{1-3} -alkoxy group, while the abovementioned unsubstituted and monosubstituted phenyl groups are additionally substituted in the 3- or 4-position

by a fluorine, chlorine or bromine atom,

by a C_{1-3} -alkoxy or C_{1-2} -alkyl-carbonyl-amino group or

by a carboxy- C_{1-3} -alkyl, aminocarbonyl- C_{1-3} -alkyl, (C_{1-2} -alkylamino)-carbonyl- C_{1-3} -alkyl, di-(C_{1-2} -alkyl)-aminocarbonyl- C_{1-3} -alkyl, (C_{1-2} -alkyl-carbonyl)-amino- C_{1-3} -alkyl or (phenyl-carbonyl)-amino- C_{1-3} -alkyl group,

while the substituents may be identical or different,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

while additionally any carboxy, amino or imino group present may be substituted by a group which can be cleaved in vivo,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

A second sub-group of compounds of general formula I deserving particular mention comprises those wherein

X denotes an oxygen atom,

R^1 denotes a hydrogen atom,

R^2 denotes a bromine atom,

R^3 denotes a phenyl group,

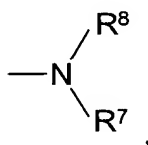
R^4 denotes a 1-(C_{1-3} -alkyl)-piperidin-4-yl group

or a phenyl group which is substituted in the 4-position

by a C₁₋₃-alkyl group terminally substituted by a C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, *N*-[ω-(di-(C₁₋₃-alkyl)-amino)-C₂₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino or *N*-(C₁₋₄-alkyloxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino group,

by a 1-(C₁₋₃-alkyl)-imidazol-2-yl or 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl group or

by a group of formula



wherein

R⁷ denotes a C₁₋₃-alkyl, C₁₋₃-alkyl-carbonyl, C₁₋₃-alkyl-sulphonyl or benzylsulphonyl group and

R⁸ denotes a ω-[di-(C₁₋₃-alkyl)-amino]-C₂₋₃-alkyl, ω-[di-(C₁₋₃-alkyl)-amino]-C₁₋₄-alkyl-carbonyl, ω-[4-(C₁₋₃-alkyl)-piperazin-1-yl]-C₁₋₃-alkyl-carbonyl or ω-{*N*-[di-(C₁₋₃-alkyl)-amino-C₂₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino}-C₁₋₃-alkyl-carbonyl group,

R⁵ denotes a hydrogen atom and

R⁶ denotes a hydrogen atom,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

A third sub-group of compounds of general formula I deserving particular mention comprises those wherein

X denotes an oxygen atom,

R¹ denotes a hydrogen atom,

R² denotes a fluorine atom,

R³ denotes a phenyl group which is optionally substituted in the 3- or 4-position by a fluorine or iodine atom or by a cyano-C₁₋₃-alkyl, amino-C₁₋₃-alkyl, C₁₋₃-alkylamino-C₁₋₃-alkyl, C₁₋₅-alkyl-carbonylamino-C₁₋₃-alkyl, C₁₋₄-alkyloxy-carbonyl-amino-C₁₋₃-alkyl, C₁₋₄-alkyloxy-C₁₋₃-alkyl-carbonyl-amino-C₁₋₃-alkyl, C₃₋₇-cycloalkyl-carbonyl-amino-C₁₋₃-alkyl, C₃₋₇-cycloalkyl-C₁₋₃-alkyl-carbonyl-amino-C₁₋₃-alkyl, *N*-(phenyl-carbonyl)-amino-C₁₋₃-alkyl, *N*-(benzyl-carbonyl)-amino-C₁₋₃-alkyl, heteroaryl-carbonyl-amino-C₁₋₃-alkyl, *N*-(C₁₋₃-alkylsulphonyl)-amino-C₁₋₃-alkyl, *N*-(phenylsulphonyl)-amino-C₁₋₃-alkyl, *N*-(benzylsulphonyl)-amino-C₁₋₃-alkyl, carboxy-C₁₋₃-alkyl, C₁₋₄-alkoxy-carbonyl-C₁₋₃-alkyl, aminocarbonyl-C₁₋₃-alkyl, C₁₋₃-alkylaminocarbonyl-C₁₋₃-alkyl, di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl-C₁₋₃-alkyl, 2-(aminocarbonyl)-C₂₋₃-alkenyl or 2-(C₁₋₃-alkyloxy-carbonyl)-C₂₋₃-alkenyl group,

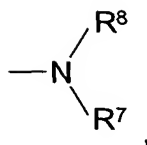
or a phenyl group trisubstituted in the 3-, 4- and 5-position by fluorine atoms,

R⁴ denotes a phenyl group which may be substituted in the 4-position

by a C₁₋₃-alkyl group terminally substituted by a pyrrolidin-1-yl, piperidin-1-yl, 4-(C₁₋₃-alkyl)-piperazin-1-yl, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₂₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino or *N*-(C₁₋₄-alkyloxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino group,

by a C₁₋₃-alkyl-sulphonyl, 1-(C₁₋₃-alkyl)-imidazol-2-yl or 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl group or

by a group of formula



wherein

R^7 denotes a C_{1-3} -alkyl, C_{1-3} -alkyl-carbonyl, C_{1-3} -alkyl-sulphonyl or benzylsulphonyl group and

R^8 denotes a C_{1-3} -alkyl, ω -[di-(C_{1-3} -alkyl)-amino]- C_{2-3} -alkyl, ω -[di-(C_{1-3} -alkyl)-amino]- C_{1-3} -alkyl-carbonyl, di-(C_{1-3} -alkyl)-amino-carbonyl- C_{1-3} -alkyl, ω -[4-(C_{1-3} -alkyl)-piperazin-1-yl]- C_{1-3} -alkyl-carbonyl or ω -{ N -[di-(C_{1-3} -alkyl)-amino- C_{2-3} -alkyl]- N -(C_{1-3} -alkyl)-amino}- C_{1-3} -alkyl-carbonyl group,

R^5 denotes a hydrogen atom and

R^6 denotes a hydrogen atom,

while the term heteroaryl group denotes a pyridinyl, furyl or thienyl group,

and unsubstituted or monosubstituted phenyl groups contained in the abovementioned definitions may additionally be substituted by a methoxy group and

the abovementioned alkyl groups include straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

A fourth sub-group of compounds of general formula I deserving particular mention comprises those wherein

X denotes an oxygen atom,

R¹ denotes a hydrogen atom,

R² denotes a cyano group,

R³ denotes a phenyl group optionally substituted by one or two methoxy groups,

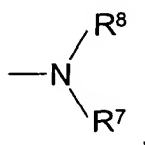
R⁴ denotes a phenyl group which is substituted in the 3- or 4-position

by a bromine atom,

by a C₁₋₃-alkyl group terminally substituted by a pyrrolidin-1-yl, piperidin-1-yl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-aminocarbonyl or *N*-(C₁₋₄-alkyloxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino group,

by a ω-di-(C₁₋₃-alkyl)-amino-C₂₋₃-alkoxy, *N*-(di-(C₁₋₃-alkyl)-amino-C₂₋₃-alkyl)-amino-carbonyl, *N*-[di-(C₁₋₃-alkyl)-amino-C₂₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino-carbonyl or 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl group or

by a group of formula



wherein

R^7 denotes a hydrogen atom or a C_{1-3} -alkyl, C_{1-3} -alkyl-carbonyl or C_{1-3} -alkylsulphonyl group and

R^8 denotes a ω -[di-(C_{1-3} -alkyl)-amino]-(C_{2-3} -alkyl), ω -[di-(C_{1-3} -alkyl)-amino]- C_{1-3} -alkyl-carbonyl, ω -(piperazin-1-yl)- C_{1-3} -alkyl-carbonyl, ω -[4-(C_{1-3} -alkyl)-piperazin-1-yl]- C_{1-3} -alkyl-carbonyl, ω -[4-(C_{1-4} -alkyloxy-carbonyl)-piperazin-1-yl]- C_{1-3} -alkyl-carbonyl, ω -[4-(C_{1-3} -alkyl)-homopiperazin-1-yl]- C_{1-3} -alkyl-carbonyl, ω -morpholino- C_{1-3} -alkyl-carbonyl or ω -{ N -[di-(C_{1-3} -alkyl)-amino]- C_{1-3} -alkyl]- N -(C_{1-3} -alkyl)-amino}- C_{1-3} -alkyl-carbonyl group,,

R^5 denotes a hydrogen atom and

R^6 denotes a hydrogen atom,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

their stereoisomers and their salts.

A fifth sub-group of compounds of general formula I deserving particular mention comprises those wherein

X denotes an oxygen or sulphur atom,

R^1 denotes a hydrogen atom,

R^2 denotes a chlorine atom,

R^3 denotes a phenyl group which is optionally monosubstituted in the 3- or 4-position

by a chlorine or iodine atom,

by a cyano, hydroxy, benzyloxy, amino or nitro group

or by an aminomethyl, acetylamino, phenylcarbonylamino-C₁₋₃-alkyl, C₁₋₃-alkylsulphonylamino-C₁₋₃-alkyl, phenylsulphonylamino-C₁₋₃-alkyl, acetylaminomethyl, imidazol-1-yl-methyl, 2-oxo-pyrrolidin-1-yl, 2-carboxy-ethyl, 2-methoxycarbonyl-ethyl, 2-aminocarbonyl-ethyl, 2-(methylaminocarbonyl)-ethyl or 2-methoxycarbonyl-ethenyl group,

or a 3-hydroxy-4-nitro-phenyl, 4-amino-3-nitrophenyl or 3,4-dimethoxyphenyl group,

R⁴ denotes a 5-(4-methyl-piperazin-1-yl-carbonyl)-pyridin-2-yl, 2-[*N*-acetyl-*N*-(ω -dimethylamino-C₂₋₃-alkyl)-amino]-pyridine-5-yl, benzo-pyrazol-6-yl, 1-methyl-2-(4-methyl-piperazin-1-yl-carbonyl)-pyrrol-4-yl, 2-(*N*-dimethylamino-ethyl-*N*-methyl-aminocarbonyl)-pyrrol-4-yl, 1-methyl-2-(*N*-dimethylamino-ethyl-*N*-methyl-aminocarbonyl)-pyrrol-4-yl, 4-(*N*-dimethylamino-methylcarbonylamino)-cyclohexyl or 4-[(*N*-dimethylamino-methylcarbonyl)-*N*-methyl-amino]cyclohexyl group or

a phenyl group which is substituted in the 3-position by a carboxy, carboxy-C₁₋₃-alkyl, C₁₋₄-alkoxy-carbonyl, C₁₋₄-alkoxy-carbonyl-C₁₋₃-alkyl, dimethylamino-C₁₋₃-alkyl or pyridin-4-yl-C₁₋₃-alkyl group or is substituted in the 4-position

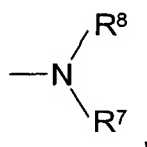
by a carboxy, ω -[di-(C₁₋₃-alkyl)-amino]-C₂₋₃-alkoxy, ethoxycarbonyl, piperidin-1-yl-carbonyl, 4-(C₁₋₄-alkyloxy-carbonyl)-piperazin-1-yl-carbonyl, *N*-(C₃₋₇-cycloalkyl)-*N*-(C₁₋₃-alkyl)-amino-carbonyl or *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-aminocarbonyl group,

by a [di-(C₁₋₃-alkyl)-amino]-C₁₋₃-alkylsulphonyl group,

by a C₁₋₃-alkyl group terminally substituted by a carboxy, C₁₋₄-alkyloxy-carbonyl, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino, *N*-(2-hydroxyethyl)-*N*-(C₁₋₃-alkyl)-amino, Di-(2-hydroxyethyl)-amino, triazolyl, *N*-(methoxyethoxyethyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-(amino-C₁₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino-carbonyl, *N*-(C₁₋₄-alkyloxy-carbonyl-amino-C₁₋₃-alkyl)-*N*-(C₁₋₃-alkyl)-amino, *N*-(C₁₋₄-alkyloxy-carbonyl)-amino or *N*-(C₁₋₄-alkyloxy-carbonyl)-*N*-(C₁₋₃-alkyl)-amino group,

by a 1-methyl-imidazol-2-yl, 5-methyl-1H-imidazol-4-yl, 1-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-imidazol-2-yl, 4-methyl-piperazin-1-yl, piperazinylcarbonyl or 4-methyl-piperazin-1-yl-carbonyl group or

by a group of formula



wherein

R⁷ denotes a hydrogen atom or a C₁₋₃-alkyl, C₁₋₃-alkyl-carbonyl, di-(C₁₋₃-alkyl)-amino-carbonyl-C₁₋₃-alkyl, benzylcarbonyl, pyridinylcarbonyl, furanylcabonyl, methoxymethylcarbonyl, C₁₋₄-alkylsulphonyl or benzylsulphonyl group or a phenylcarbonyl group optionally substituted in the phenyl moiety by one or two methoxy groups and

R⁸ denotes a C₁₋₃-alkyl, ω-[di-(C₁₋₃-alkyl)-amino]-C₂₋₃-alkyl, ω-[*N*-benzyl-*N*-(C₁₋₃-alkyl)-amino]-C₂₋₃-alkyl, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-amino-carbonyl, (pyridinyl-C₁₋₃-alkyl)-amino-carbonyl, 1-(C₁₋₃-alkyl)-piperidin-4-yl-amino-carbonyl-, 1-(C₁₋₃-alkyl)-piperidin-4-yl-oxy-carbonyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-amino-

carbonyl, 4-(C₁₋₃-alkyl)-piperazin-1-yl-carbonyl or di-(C₁₋₃-alkyl)-aminocarbonyl-C₁₋₃-alkyl group or

a C₁₋₄-alkyl-carbonyl group terminally substituted by a hydroxy, amino, di-(C₁₋₃-alkyl)-amino, *N*-benzyl-*N*-(C₁₋₃-alkyl)-amino, di-(2-hydroxyethyl)-amino, acetylamino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-(C₁₋₃-alkyl)-amino, imidazol-1-yl, piperazin-1-yl, 4-(C₁₋₃-alkyl)-piperazin-1-yl, 4-benzyl-piperazin-1-yl, 4-(C₁₋₄-alkyloxy-carbonyl)-piperazin-1-yl, 4-(C₁₋₃-alkyl)-homopiperazin-1-yl, morpholin-4-yl, pyrrolidin-1-yl, piperidin-1-yl, 1-(C₁₋₃-alkyl)-piperidin-4-yl or phthalimido group,

R⁵ denotes a hydrogen atom and

R⁶ denotes a hydrogen atom or a nitro group,

while the unsubstituted or monosubstituted phenyl groups mentioned in the above definitions may additionally be substituted by a methoxy or a cyano group or by two methyl groups,

the abovementioned alkyl groups including straight-chain and branched alkyl groups, wherein additionally one to 3 hydrogen atoms may be replaced by fluorine atoms,

the tautomers, enantiomers, diastereomers, mixtures thereof and the salts thereof.

Additional sub-groups of compounds of general formula I, of preferred, particularly preferred and most particularly preferred compounds of formula I and the sub-groups thereof which deserve special mention are those wherein X, R¹, R³, R⁴, R⁵ and R⁶ are each as hereinbefore defined and R² denotes fluorine.

Additional sub-groups of compounds of general formula I, of preferred, particularly preferred and most particularly preferred compounds of formula I and the sub-groups thereof which deserve special mention are those wherein X, R¹, R³, R⁴, R⁵ and R⁶ are each as hereinbefore defined and R² denotes chlorine.

Additional sub-groups of compounds of general formula I, of preferred, particularly preferred and most particularly preferred compounds of formula I and the sub-groups thereof which deserve special mention are those wherein X, R¹, R³, R⁴, R⁵ and R⁶ are each as hereinbefore defined and R² denotes bromine.

Additional sub-groups of compounds of general formula I, of preferred, particularly preferred and most particularly preferred compounds of formula I and the sub-groups thereof which deserve special mention are those wherein X, R¹, R³, R⁴, R⁵ and R⁶ are each as hereinbefore defined and R² denotes cyano.

The following compounds of general formula I are particularly preferred:

- (a) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone,
- (b) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone,
- (c) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone,
- (d) 3-Z-[1-(4-(N-(4-ethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone,
- (e) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone,

- (f) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone,
- (g) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone,
- (h) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone,
- (i) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone,
- (j) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone,
- (k) 3-Z-[1-(4-(N-(4-ethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone,
- (l) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone,
- (m) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone,
- (n) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone,
- (o) 3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone,
- (p) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone,

(q) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone,

(r) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-carbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone,

(s) 3-Z-[1-(4-(N-(piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone,

(t) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone,

(u) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone,

(v) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone,

(w) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone,

(x) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone,

(y) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone,

(z) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-dimethylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone,

(aa) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-dimethylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone,

(ab) 3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone,

(ac) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-acetylamino-phenyl)-methylene]-6-chloro-2-indolinone,

(ad) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-acetylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone,

(ae) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-acetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone,

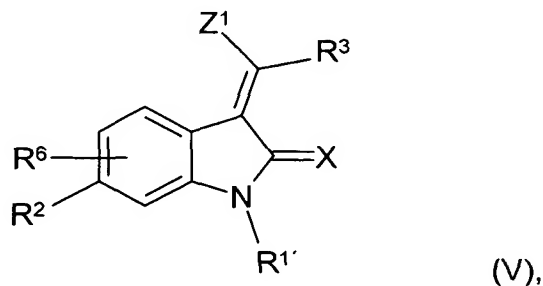
(af) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-benzoylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone and

(ag) 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-acetylamino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

and the salts thereof.

According to the invention the new compounds may be obtained for example by the following methods known in principle from the literature:

a. reacting a compound of general formula



wherein

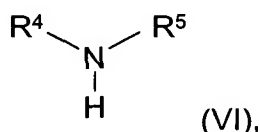
the groups Z^1 and R^3 may optionally change positions,

X, R², R³ and R⁶ are as hereinbefore defined,

R^{1a} has the meanings given for R¹ hereinbefore or denotes a protective group for the nitrogen atom of the lactam group, while R¹ may also denote a bond to a solid phase, optionally via a spacer,

and Z¹ denotes a halogen atom, a hydroxy, alkoxy or aryl-alkoxy group, e.g. a chlorine or bromine atom, a methoxy, ethoxy or benzyloxy group,

with an amine of general formula



wherein

R⁴ and R⁵ are as hereinbefore defined,

and if necessary subsequently cleaving any protective group used for the nitrogen atom of the lactam group or from a solid phase.

The protective group for the nitrogen atom of the lactam group may be for example an acetyl, benzoyl, ethoxycarbonyl, tert.butyloxycarbonyl or benzyloxycarbonyl group and

the solid phase may be a resin such as a 4-(2',4'-dimethoxyphenylamino-methyl)-phenoxy resin, while the bond is conveniently made via the amino group, or a p-benzyloxybenzylalcohol resin, while the bond is conveniently made via an intermediate member such as a 2,5-dimethoxy-4-hydroxy-benzyl derivative.

The reaction is expediently carried out in a solvent such as dimethylformamide, toluene, acetonitrile, tetrahydrofuran, dimethylsulphoxide, methylene chloride or mixtures thereof, optionally in the presence of an inert base such as triethylamine, N-ethyl-diisopropylamine or sodium hydrogen carbonate at temperatures between 20 and 175°C, while any protective group used may simultaneously be cleaved as a result of transamidation.

If Z^1 in a compound of general formula V denotes a halogen atom, then the reaction is preferably carried out in the presence of an inert base at temperatures between 20 and 120°C.

If Z^1 in a compound of general formula V denotes a hydroxy, alkoxy or arylalkoxy group, then the reaction is preferably carried out at temperatures between 20 and 200°C.

Any protective group used is optionally subsequently conveniently cleaved either by hydrolysis in an aqueous or alcoholic solvent, e.g. in methanol/water, ethanol/water, isopropanol/water, tetrahydrofuran/water, dioxane/water, dimethylformamide/water, methanol or ethanol in the presence of an alkali metal base such as lithium hydroxide, sodium hydroxide or potassium hydroxide at temperatures between 0 and 100°C, preferably at temperatures between 10 and 50°C,

or advantageously by transamidation with an organic base such as ammonia, butylamine, dimethylamine or piperidine in a solvent such as methanol, ethanol, dimethylformamide and mixtures thereof or in an excess of the amine used at temperatures between 0 and 100°C, preferably at temperatures between 10 and 50°C.

The cleaving of any solid phase used is preferably carried out using trifluoroacetic acid and water at temperatures between 0 and 35°C, preferably at ambient temperature.

b. In order to prepare a compound of general formula I wherein R^4 contains the group R^8 , where

R^8 denotes a C_{1-4} -alkyl-carbonyl group terminally substituted by a hydroxy, C_{1-3} -alkyoxyl group, amino, (C_{1-3} -alkyl)-amino, di- (C_{1-3} -alkyl)-amino, (ω -hydroxy- C_{2-3} -alkyl)-amino, di-(ω -hydroxy- C_{2-3} -alkyl)-amino, (ω -alkoxy- C_{2-3} -alkyl)-amino, di-(ω -alkoxy- C_{2-3} -

alkyl)-amino, C₁₋₃-alkyl-carbonyl-amino, *N*-benzyl-*N*-C₁₋₃-alkyl-amino, *N*-[di-(C₁₋₃-alkyl)-amino-C₁₋₃-alkyl]-*N*-C₁₋₃-alkyl-amino, 1-(C₁₋₃-alkyl)-piperidin-4-yl- group or by a 5- to 7-membered cycloalkyleneimino group,

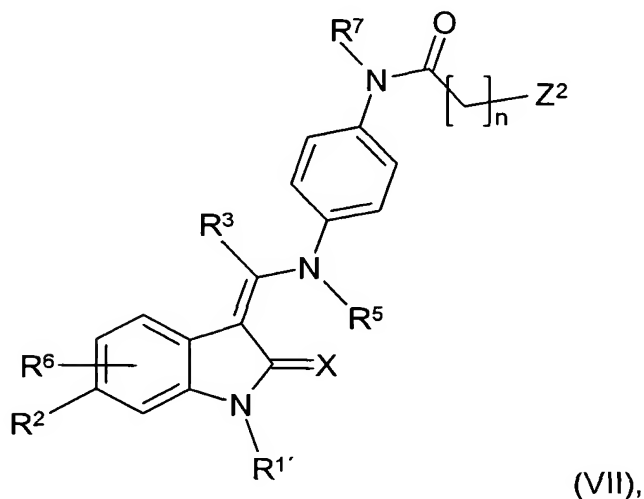
while the cycloalkylene group may be substituted by a C₁₋₃-alkyl group and/or

one or two methylene groups linked to the imino group may be replaced by a carbonyl group and/or

the methylene group in the 4 position of a 6- or 7-membered cycloalkylimino group may be replaced by a –NH, –N(C₁₋₃-alkyl), –N(benzyl), –N(C₁₋₄-alkoxy-carbonyl) or –O- and/or

a phenyl ring may be fused on via two adjacent carbon atoms of the cycloalkyleneimino group:

reacting a compound of general formula



wherein

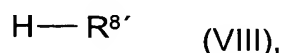
R², R³, R⁵, R⁶, R⁷ and X are as hereinbefore defined,

$R^{1'}$ has the meanings given for R^1 hereinbefore or denotes a protective group for the nitrogen atom of the lactam group, while $R^{1'}$ may also denote a bond to a solid phase optionally formed via a spacer,

n denotes the number 1, 2, 3 or 4 and

Z^2 denotes a leaving group, for example a halogen atom or an alkyl or arylsulphonyloxy group such as the chlorine, bromine or iodine atom or the methylsulphonyloxy, ethylsulphonyloxy, p-toluenesulphonyloxy, or trifluoromethanesulphonyloxy group,

with a hydroxide base such as sodium or potassium hydroxide or a compound of general formula



wherein

$R^{8'}$ denotes a C_{1-3} -alkyloxy, amino, $(C_{1-3}$ -alkyl)-amino, di- $(C_{1-3}$ -alkyl)-amino, $(\omega$ -hydroxy- C_{2-3} -alkyl)-amino, di- $(\omega$ -hydroxy- C_{2-3} -alkyl)-amino, $(\omega$ -alkoxy- C_{2-3} -alkyl)-amino, di- $(\omega$ -alkoxy- C_{2-3} -alkyl)-amino, C_{1-3} -alkyl-carbonyl-amino, *N*-benzyl-*N*- C_{1-3} -alkyl-amino, *N*-[di- $(C_{1-3}$ -alkyl)-amino- C_{1-3} -alkyl]-*N*- C_{1-3} -alkyl-amino or a 5- to 7-membered cycloalkyleneimino group,

while the cycloalkylene group may be substituted by a C_{1-3} -alkyl group and/or

one or two methylene groups linked to the imino group may be replaced by a carbonyl group and/or

the methylene group in the 4 position of a 6- or 7-membered cycloalkylimino group may be replaced by a -NH, -N(C_{1-3} -alkyl), -N(benzyl), -N(C_{1-4} -alkoxy-carbonyl) or -O- and/or

a phenyl ring may be fused on via two adjacent carbon atoms of the cycloalkyleneimino group,

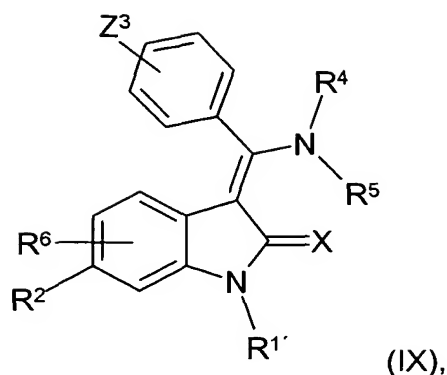
and if necessary subsequently cleaving any protective group used for the nitrogen atom of the lactam group or from a solid phase.

The reaction is expediently carried out in a solvent such as methylene chloride, tetrahydrofuran, 1,4-dioxane, toluene, acetonitrile, dimethylsulphoxide, dimethylformamide, dimethylacetamide, N-methylpyrrolidone or mixtures thereof, optionally with the addition of water as cosolvent and/or with the addition of an inert auxiliary base, for example sodium hydrogen carbonate, pyridine, 2,4,6-trimethylpyridine, quinoline, triethylamine, N-ethyldiisopropylamine, N-ethyl-dicyclohexylamine, 1,4-diazabicyclo[2,2,2]octane or 1,8-diazabicyclo[5,4,0]undec-7-ene, at temperatures between -50°C and +100°C, preferably between -10°C and +50°C, while any protective group used may simultaneously be cleaved as a result of transamidation.

If it is necessary to cleave a protective group used for the nitrogen atom of the lactam group or to cleave from a solid phase the same method is used as described under (a) above.

c. In order to prepare a compound of general formula I wherein R^3 denotes a phenyl or naphthyl group substituted by a carboxy- C_{2-3} -alkenyl, amino-carbonyl- C_{2-3} -alkenyl, (C_{1-3} -alkylamino)-carbonyl- C_{2-3} -alkenyl, di-(C_{1-3} -alkylamino)-carbonyl- C_{2-3} -alkenyl or C_{1-4} -alkoxy-carbonyl- C_{2-3} -alkenyl group,

reacting a compound of general formula



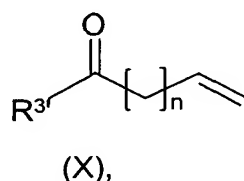
wherein

R^2 , R^4 , R^5 , R^6 and X are as hereinbefore defined,

$R^{1'}$ has the meanings given for R^1 hereinbefore or denotes a protective group for the nitrogen atom of the lactam group, while $R^{1'}$ may also denote a bond to a solid phase optionally formed via a spacer, and

Z^3 denotes a leaving group, for example a halogen atom or an alkyl or arylsulphonyloxy group such as the chlorine, bromine or iodine atom or the methylsulphonyloxy, ethylsulphonyloxy, p-toluenesulphonyloxy or trifluoromethanesulphonyloxy group,

with an alkene of general formula



wherein

$R^{3'}$ denotes an amino, (C₁₋₃-alkylamino), di-(C₁₋₃-alkylamino) or C₁₋₄-alkoxy-group and n is the number 0 or 1.

The reaction is conveniently carried out under palladium catalysis, for example with palladium(II)-acetate, palladium(II)-chloride, bis-(triphenylphosphine)-palladium(II)-acetate, bis-(triphenylphosphine)-palladium(II)-chloride, palladium/activated charcoal, bis-[1,2-bis-(diphenylphosphino)-ethane]-palladium(0), dichloro-(1,2-bis-(diphenylphosphino)-ethane)-palladium(II), tetrakis(triphenylphosphine)-palladium(0), tris-(dibenzylideneacetone)-dipalladium(0), 1,1'-bis-

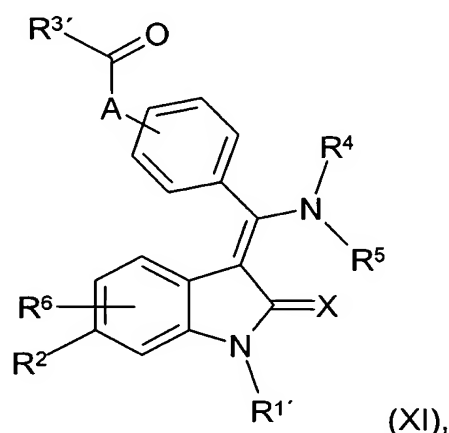
(diphenylphosphino)-ferrocene-dichloro-palladium(II) or tris-(dibenzylideneacetone)-dipalladium(0)-chloroform adduct in the presence of a base such as triethylamine, diisopropyl-ethylamine, lithium carbonate, potassium carbonate, sodium carbonate, caesium carbonate and a ligand such as triphenylphosphine, tri-ortho-tolyl-phosphine or tri-(tert.butyl)-phosphine in solvents such as acetonitrile, N-methyl-pyrrolidinone, dioxane or dimethylformamide and mixtures thereof.

If it is necessary to cleave a protective group used for the nitrogen atom of the lactam group or to cleave from a solid phase the same method is used as described under (a) above.

d. In order to prepare a compound of general formula I wherein R^3 denotes a phenyl or naphthyl group substituted by

carboxy- C_{1-3} -alkyl, C_{1-4} -alkoxy-carbonyl- C_{1-3} -alkyl, aminocarbonyl- C_{1-3} -alkyl, (C_{1-3} -alkylamino)-carbonyl- C_{1-3} -alkyl, di-(C_{1-3} -alkyl)-amino-carbonyl- C_{1-3} -alkyl or 4-(C_{1-3} -alkyl)-piperazin-1-yl-carbonyl-(C_{1-3} -alkyl) groups,

hydrogenating a compound of general formula



wherein

R^2 , R^4 , R^5 , R^6 and X are as hereinbefore defined,

$R^{1'}$ has the meanings given for R^1 hereinbefore or denotes a protective group for the nitrogen atom of the lactam group, while $R^{1'}$ may also denote a bond to a solid phase optionally formed via a spacer,

A denotes a C_{2-3} -alkenyl group and

$R^{3'}$ denotes a hydroxy, C_{1-4} -alkoxy, amino, (C_{1-3} -alkylamino), di-(C_{1-3} -alkyl)-amino or 4-(C_{1-3} -alkyl)-piperazin-1-yl group.

The hydrogenation is preferably carried out by catalytic hydrogenation with hydrogen in the presence of a catalyst such as palladium/charcoal or platinum in a solvent such as methanol, ethanol, ethyl acetate, dimethylformamide, dimethylformamide/acetone or glacial acetic acid optionally with the addition of an acid such as hydrochloric acid at temperatures between 0 and 50°C, but preferably at ambient temperature, and under a hydrogen pressure of 1 to 7 bar, but preferably from 3 to 5 bar.

If it is necessary to cleave a protective group used for the nitrogen atom of the lactam group or to cleave from a solid phase the same method is used as described under (a) above.

If according to the invention a compound of general formula I is obtained which contains an alkoxycarbonyl group, this may be converted by hydrolysis into a corresponding carboxy compound, or

if a compound of general formula I is obtained which contains an amino or alkylamino group, this may be converted by reductive alkylation into a corresponding alkylamino or dialkylamino compound, or

if a compound of general formula I is obtained which contains an amino or alkylamino group, this may be converted by acylation or sulphonation into a corresponding acyl or sulphonyl compound, or

if a compound of general formula I is obtained which contains a carboxy group, this may be converted by esterification or amidation into a corresponding ester or aminocarbonyl compound, or

if a compound of general formula I is obtained which contains a cyclo-alkyleneimino group in which a methylene group is replaced by a sulphur atom, this may be converted by oxidation into a corresponding sulphinyl or sulphonyl compound, or

if a compound of general formula I is obtained which contains a nitro group, this may be converted by reduction into a corresponding amino compound, or

if a compound of general formula I is obtained which contains a cyano group, this may be converted by reduction into a corresponding aminomethyl compound, or

if a compound of general formula I is obtained which contains an arylalkyloxy group, this may be converted with acid into a corresponding hydroxy compound, or

if a compound of general formula I is obtained which contains an alkoxycarbonyl group, this may be converted by saponification into a corresponding carboxy compound, or

if a compound of general formula I is obtained wherein R_4 denotes a phenyl group substituted by an amino, alkylamino, aminoalkyl or N-alkyl-amino group, this may be converted by reacting with a corresponding cyanate, isocyanate or carbamoyl halide into a corresponding urea compound of general formula I or

if a compound of general formula I is obtained which contains a carbonyl group, this may be converted by reacting with phosphorus pentasulphide into a corresponding thiocarbonyl compound, or

if a compound of general formula I is obtained wherein R_4 denotes a phenyl group substituted by an amino, alkylamino, aminoalkyl or N-alkyl-amino group, this may subsequently be converted by reacting with a corresponding

compound which transfers the amidino group or by reacting with a corresponding nitrile into a corresponding guanidino compound of general formula I.

The subsequent hydrolysis is preferably carried out in an aqueous solvent, e.g. in water, methanol/water, ethanol/water, isopropanol/water, tetrahydrofuran/water or dioxane/water, in the presence of an acid such as trifluoroacetic acid, hydrochloric acid or sulphuric acid or in the presence of an alkali metal base such as lithium hydroxide, sodium hydroxide or potassium hydroxide at temperatures between 0 and 100°C, preferably at temperatures between 10 and 50°C.

The subsequent reductive alkylation is preferably carried out in a suitable solvent such as methanol, methanol/water, methanol/water/ammonia, ethanol, ether, tetrahydrofuran, dioxane or dimethylformamide, optionally with the addition of an acid such as hydrochloric acid in the presence of catalytically activated hydrogen, e.g. hydrogen in the presence of Raney nickel, platinum or palladium/charcoal, or in the presence of a metal hydride such as sodium borohydride, lithium borohydride, sodium cyanoborohydride or lithium aluminium hydride at temperatures between 0 and 100°C, preferably at temperatures between 20 and 80°C.

The subsequent acylation or sulphonylation is conveniently carried out with the corresponding free acid or a corresponding reactive compound such as the anhydride, ester, imidazolide or halide thereof, preferably in a solvent such as methylene chloride, diethyl ether, tetrahydrofuran, toluene, dioxane, acetonitrile, dimethylsulphoxide or dimethylformamide, optionally in the presence of an inorganic or a tertiary organic base at temperatures between -20 and 200°C, preferably at temperatures between 20°C and the boiling temperature of the solvent used. The reaction with the free acid may optionally take place in the presence of an acid-activating agent or a dehydrating agent, e.g. in the presence of isobutyl chloroformate, tetraethyl orthocarbonate, trimethyl orthoacetate, 2,2-dimethoxypropane, tetramethoxysilane, thionyl chloride, trimethylchlorosilane, phosphorus

trichloride, phosphorus pentoxide, N,N'-dicyclohexylcarbodiimide, N,N'-dicyclohexylcarbodiimide/N-hydroxysuccinimide, N,N'-dicyclohexylcarbodiimide/1-hydroxy-benzotriazole, 2-(1H-benzotriazol-1-yl)-1,1,3,3-tetramethyluronium-tetrafluoroborate, 2-(1H-benzotriazol-1-yl)-1,1,3,3-tetramethyluronium-tetrafluoroborate/1-hydroxy-benzotriazole, N,N'-carbonyldiimidazole or triphenylphosphine/carbon tetrachloride, and optionally with the addition of a base such as pyridine, 4-dimethylamino-pyridine, N-methyl-morpholine or triethylamine, conveniently at temperatures between 0 and 150°C, preferably at temperatures between 0 and 100°C. The reaction with a corresponding reactive compound may optionally take place in the presence of a tertiary organic base such as triethylamine, N-ethyl-diisopropylamine, N-methyl-morpholine or pyridine or using an anhydride in the presence of the corresponding acid at temperatures between 0 and 150°C, preferably at temperatures between 50 and 100°C.

The subsequent esterification or amidation is conveniently carried out by reacting a corresponding reactive carboxylic acid derivative with a corresponding alcohol or amine as hereinbefore described.

The esterification or amidation is preferably carried out in a solvent such as methylene chloride, diethyl ether, tetrahydrofuran, toluene, dioxane, acetonitrile, dimethylsulphoxide or dimethylformamide, optionally in the presence of an inorganic or a tertiary organic base, preferably at temperatures between 20°C and the boiling temperature of the solvent used. The reaction is carried out with a corresponding acid, preferably in the presence of a dehydrating agent, e.g. in the presence of isobutyl chloroformate, tetraethyl orthocarbonate, trimethyl orthoacetate, 2,2-dimethoxypropane, tetramethoxysilane, thionyl chloride, trimethylchlorosilane, phosphorus trichloride, phosphorus pentoxide, N,N'-dicyclohexylcarbodiimide, N,N'-dicyclohexylcarbodiimide/N-hydroxysuccinimide, N,N'-dicyclohexylcarbodiimide/1-hydroxy-benzotriazole, 2-(1H-benzotriazol-1-yl)-1,1,3,3-tetramethyluronium-tetrafluoroborate, 2-(1H-benzotriazol-1-yl)-1,1,3,3-tetramethyluronium-tetrafluoroborate/1-hydroxy-benzotriazole, N,N'-carbonyldiimidazole or triphenylphosphine/carbon tetrachloride, and

optionally with the addition of a base such as pyridine, 4-dimethylaminopyridine, N-methyl-morpholine or triethylamine, conveniently at temperatures between 0 and 150°C, preferably at temperatures between 0 and 100°C, and the acylation with a corresponding reactive compound such as the anhydride, ester, imidazolid or halide thereof may optionally be carried out in the presence of a tertiary organic base such as triethylamine, N-ethyl-diisopropylamine or N-methyl-morpholine at temperatures between 0 and 150°C, preferably at temperatures between 50 and 100°C.

The subsequent oxidation of the sulphur atom is preferably carried out in a solvent or mixture of solvents, e.g. in water, water/pyridine, acetone, methylene chloride, acetic acid, acetic acid/acetic anhydride, dilute sulphuric acid or trifluoroacetic acid, conveniently at temperatures between -80 and 100°C, depending on the oxidising agent used.

In order to prepare a corresponding sulphinyl compound of general formula I the oxidation is conveniently carried out with one equivalent of the oxidising agent used, e.g. with hydrogen peroxide in glacial acetic acid, trifluoroacetic acid or formic acid at 0 to 20°C or in acetone at 0 to 60°C, with a peracid such as performic acid in glacial acetic acid or trifluoroacetic acid at 0 to 50°C or with m-chloroperbenzoic acid in methylene chloride, chloroform or dioxane at -20 to 80°C, with sodium metaperiodate in aqueous methanol or ethanol at -15 to 25°C, with bromine in glacial acetic acid or aqueous acetic acid optionally in the presence of a weak base such as sodium acetate, with N-bromosuccinimide in ethanol, with tert.butylhypochlorite in methanol at -80 to -30°C, with iodobenzodichloride in aqueous pyridine at 0 to 50°C, with nitric acid in glacial acetic acid at 0 to 20°C, with chromic acid in glacial acetic acid or in acetone at 0 to 20°C and with sulphurylchloride in methylene chloride at -70°C, and the thioether-chlorine complex thus obtained is conveniently hydrolysed with aqueous ethanol.

In order to prepare a sulphonyl compound of general formula I oxidation is carried out starting from a corresponding sulphinyl compound, conveniently with one or more equivalents of the oxidising agent used, or starting from a

corresponding mercapto compound, conveniently with two or more equivalents of the oxidising agent used, e.g. with hydrogen peroxide in glacial acetic acid/acetic anhydride, trifluoroacetic acid or in formic acid at 20 to 100°C or in acetone at 0 to 60°C, with a peracid such as performic acid or m-chloroperbenzoic acid in glacial acetic acid, trifluoroacetic acid, methylene chloride or chloroform at temperatures between 0 and 60°C, with nitric acid in glacial acetic acid at 0 to 20°C, with chromic acid, sodium periodate or potassium permanganate in acetic acid, water/sulphuric acid or in acetone at 0 to 20°C.

The subsequent reduction of a nitro group is preferably carried out by hydrogenolysis, e.g. with hydrogen in the presence of a catalyst such as palladium/charcoal or Raney nickel in a solvent such as methanol, ethanol, ethyl acetate, dimethylformamide, dimethylformamide/acetone or glacial acetic acid, optionally with the addition of an acid such as hydrochloric acid or glacial acetic acid at temperatures between 0 and 50°C, but preferably at ambient temperature, and under a hydrogen pressure of 1 to 7 bar, but preferably 3 to 5 bar.

The subsequent hydrogenation of a cyano group is preferably carried out by hydrogenolysis, e.g. with hydrogen in the presence of a catalyst such as palladium/charcoal or Raney nickel in a solvent such as methanol, ethanol, ethyl acetate, methylene chloride, dimethylformamide, dimethylformamide/acetone or glacial acetic acid, optionally with the addition of an acid such as hydrochloric acid or glacial acetic acid at temperatures between 0 and 50°C, but preferably at ambient temperature, and under a hydrogen pressure of 1 to 7 bar, but preferably from 3 to 5 bar.

Expediently, a corresponding urea compound of general formula I is subsequently prepared with an inorganic cyanate or a corresponding isocyanate or carbamoylchloride, preferably in a solvent such as dimethylformamide and optionally in the presence of a tertiary organic base such as triethylamine at temperatures between 0 and 50°C, preferably at ambient temperature.

Expediently, a corresponding thiocarbonyl compound of general formula I is subsequently prepared with phosphorus pentasulphide or (p-methoxy-phenyl)-thionophosphine-sulphide dimer (Lawesson's reagent) preferably in a solvent such as pyridine or toluene at temperatures between 80 and 120°C, preferably at 120°C.

Expediently, a corresponding guanidino compound of general formula I is subsequently prepared by reacting with a compound which transfers the amidino group, such as 3,5-dimethylpyrazole-1-carboxylic acid amidine, preferably in a solvent such as dimethylformamide and optionally in the presence of a tertiary organic base such as triethylamine at temperatures between 0 and 50°C, preferably at ambient temperature.

In the reactions described hereinbefore, any reactive groups present such as carboxy, hydroxy, amino, alkylamino or imino groups may be protected during the reaction by conventional protecting groups which are cleaved again after the reaction.

For example, a protecting group for a carboxy group may be a trimethylsilyl, methyl, ethyl, tert.butyl, benzyl or tetrahydropyranyl group and

protecting groups for a hydroxy, amino, alkylamino or imino group may be an acetyl, trifluoroacetyl, benzoyl, ethoxycarbonyl, tert.butoxycarbonyl, benzyloxycarbonyl, benzyl, methoxybenzyl or 2,4-dimethoxybenzyl group and additionally, for the amino group, a phthalyl group.

Any protecting group used is optionally subsequently cleaved for example by hydrolysis in an aqueous solvent, e.g. in water, isopropanol/water, tetrahydrofuran/water or dioxan/water, in the presence of an acid such as trifluoroacetic acid, hydrochloric acid or sulphuric acid or in the presence of an alkali metal base such as lithium hydroxide, sodium hydroxide or potassium hydroxide at temperatures between 0 and 100°C, preferably at temperatures between 10 and 50°C.

However, a benzyl, methoxybenzyl or benzyloxycarbonyl group is cleaved, for example, hydrogenolytically, e.g. with hydrogen in the presence of a catalyst such as palladium/charcoal in a solvent such as methanol, ethanol, ethyl acetate, dimethylformamide, dimethylformamide/acetone or glacial acetic acid, optionally with the addition of an acid such as hydrochloric acid or glacial acetic acid at temperatures between 0 and 50°C, but preferably at ambient temperature, and at a hydrogen pressure of 1 to 7 bar, but preferably 3 to 5 bar.

A methoxybenzyl group may also be cleaved in the presence of an oxidising agent such as cerium(IV)ammonium nitrate in a solvent such as methylene chloride, acetonitrile or acetonitrile/water at temperatures of between 0 and 50°C, but preferably at ambient temperature.

A 2,4-dimethoxybenzyl group, however, is preferably cleaved in trifluoroacetic acid in the presence of anisole.

A tert.butyl or tert.butyloxycarbonyl group is preferably cleaved by treating with an acid such as trifluoroacetic acid or hydrochloric acid, optionally using a solvent such as methylene chloride, dioxane, ethyl acetate or ether.

A phthalyl group is preferably cleaved in the presence of hydrazine or a primary amine such as methylamine, ethylamine or n-butylamine in a solvent such as methanol, ethanol, isopropanol, toluene/water or dioxane at temperatures between 20 and 50°C.

Moreover, chiral compounds of general formula I obtained may be resolved into their enantiomers and/or diastereomers.

Thus, for example, the compounds of general formula I obtained which occur as racemates may be separated by methods known per se (cf. Allinger N. L. and Eliel E. L. in "Topics in Stereochemistry", Vol. 6, Wiley Interscience, 1971) into their optical antipodes and compounds of general formula I with at least 2

asymmetric carbon atoms may be resolved into their diastereomers on the basis of their physical-chemical differences using methods known per se, e.g. by chromatography and/or fractional crystallisation, and, if these compounds are obtained in racemic form, they may subsequently be resolved into the enantiomers as mentioned above.

The enantiomers are preferably separated by column separation on chiral phases or by recrystallisation from an optically active solvent or by reacting with an optically active substance which forms salts or derivatives such as e.g. esters or amides with the racemic compound, particularly acids and the activated derivatives or alcohols thereof, and separating the diastereomeric mixture of salts or derivatives thus obtained, e.g. on the basis of their differences in solubility, whilst the free antipodes may be released from the pure diastereomeric salts or derivatives by the action of suitable agents. Optically active acids in common use are e.g. the D- and L-forms of tartaric acid or dibenzoyltartaric acid, di-o-tolytartaric acid, malic acid, mandelic acid, camphorsulphonic acid, glutamic acid, N-acetylglutamic acid, aspartic acid, N-acetylaspartic acid or quinic acid. An optically active alcohol may be for example (+) or (-)-menthol and an optically active acyl group in amides, for example, may be a (+)-or (-)-menthyloxycarbonyl group.

Furthermore, the compounds of formula I obtained may be converted into the salts thereof, particularly for pharmaceutical use into the physiologically acceptable salts with inorganic or organic acids. Acids which may be used for this purpose include for example hydrochloric acid, hydrobromic acid, sulphuric acid, phosphoric acid, fumaric acid, succinic acid, lactic acid, citric acid, tartaric acid, maleic acid, methanesulphonic acid, ethanesulphonic acid, para-toluenesulphonic acid, phenylsulphonic acid or L-(+)-mandelic acid.

Moreover, if the new compounds of formula I thus obtained contain a carboxy group, they may subsequently, if desired, be converted into the salts thereof with inorganic or organic bases, particularly for pharmaceutical use into the physiologically acceptable salts thereof. Suitable bases for this purpose

include for example sodium hydroxide, potassium hydroxide, cyclohexylamine, ethanolamine, diethanolamine and triethanolamine.

For compounds of general formula I which contain 2 or more acidic or basic groups, salts with 2 or more inorganic or organic bases or acids may also be used (so-called disalts etc.).

The compounds of general formulae V to XI used as starting products are known from the literature in some cases or may be obtained by methods known from the literature or may be obtained by methods described hereinbefore and in the Examples. For example, the compounds of general formula IX are described in German Patent Application 198 44 003.

As already mentioned hereinbefore, the new compounds of general formula I have valuable pharmacological properties, particularly an inhibiting effect on various kinases, especially receptor tyrosine kinases such as VEGFR1, VEGFR2, VEGFR3, PDGFR α , PDGFR β , FGFR1, FGFR3, EGFR, HER2, c-Kit, IGF1R and HGFR, Flt-3, as well as on the proliferation of cultivated human cells, particularly endothelial cells, e.g. in angiogenesis, but also on the proliferation of other cells, particularly tumour cells.

The biological properties of the new compounds were tested by the following standard method as described:

Human umbilical endothelial cells (HUVEC) were cultivated in IMDM (Gibco BRL), supplemented with 10 % foetal calf serum (FBS) (Sigma), 50 μ M of β -mercaptoethanol (Fluka), standard antibiotics, 15 μ g/ml of endothelial cell growth factor (ECGS, Collaborative Biomedical Products) and 100 μ g/ml of heparin (Sigma) on gelatine-coated culture dishes (0.2 % gelatine, Sigma) at 37°C, under 5 % CO₂ in a water-saturated atmosphere.

In order to investigate the inhibitory activity of the compounds according to the invention the cells were "starved" for 16 hours, i.e. kept in culture medium without growth factors (ECGS + heparin). The cells were detached from the culture dishes using trypsin/EDTA and washed once in serum-containing medium. Then they were seeded out in amounts of 2.5×10^3 cells per well.

The proliferation of the cells was stimulated with 5 ng/ml of VEGF₁₆₅ (vascular endothelial growth factor; H. Weich, GBF Braunschweig) and 10 µg/ml of heparin. As a control, 6 wells in each dish were not stimulated.

The compounds according to the invention were dissolved in 100% dimethylsulphoxide and added to the cultures in various dilutions as triple measurements, the maximum dimethylsulphoxide concentration being 0.3 %.

The cells were incubated for 76 hours at 37°C, then for a further 16 hours ³H-thymidine (0.1 µ Ci/well, Amersham) was added in order to determine the DNA synthesis. Then the radioactively labelled cells were immobilised on filter mats and the radioactivity incorporated was measured in a β-counter. In order to determine the inhibitory activity of the compounds according to the invention the mean value of the non-stimulated cells was subtracted from the mean value of the factor-stimulated cells (in the presence or absence of the compounds according to the invention).

The relative cell proliferation was calculated as a percentage of the control (HUVEC without inhibitor) and the concentration of active substance which inhibits the proliferation of the cells by 50 % (IC₅₀) was determined.

The compounds of general formula I according to the invention have an IC₅₀ of between 50 µM and 1 nM.

In view of their inhibitory effect on the proliferation of cells, particularly endothelial and tumour cells, the compounds of general formula I are suitable

for treating diseases in which the proliferation of cells, particularly endothelial cells, plays a part.

Thus, for example, the proliferation of endothelial cells and the concomitant neovascularisation constitute a crucial stage in tumour progression (Folkman J. et al., *Nature* 339, 58-61, (1989); Hanahan D. and Folkman J., *Cell* 86, 353-365, (1996)). Furthermore, the proliferation of endothelial cells is also important in haemangiomas, in metastatisation, rheumatoid arthritis, psoriasis and ocular neovascularisation (Folkman J., *Nature Med.* 1, 27-31, (1995)). The therapeutic usefulness of inhibitors of endothelial cell proliferation was demonstrated in the animal model for example by O'Reilly et al. and Parangi et al. (O'Reilly M.S. et al., *Cell* 88, 277-285, (1997); Parangi S. et al., *Proc Natl Acad Sci USA* 93, 2002-2007, (1996)).

The compounds of general formula I, their tautomers, their stereoisomers or the physiologically acceptable salts thereof are thus suitable, for example, for treating tumours (e.g. plate epithelial carcinoma, astrocytoma, Kaposi's sarcoma, glioblastoma, lung cancer, bladder cancer, head and neck carcinoma, melanoma, ovarian cancer, prostate cancer, breast cancer, small-cell lung cancer, glioma, colorectal cancer, urogenital cancer and gastrointestinal cancer as well as haematological cancers such as multiple myeloma), psoriasis, arthritis (e.g. rheumatoid arthritis), haemangioma, angiofibroma, eye diseases (e.g. diabetic retinopathy), neovascular glaucoma, kidney diseases (e.g. glomerulonephritis), diabetic nephropathy, malignant nephrosclerosis, thrombotic microangiopathic syndrome, transplant rejections and glomerulopathy, fibrotic diseases (e.g. cirrhosis of the liver), mesangial cell proliferative diseases, arteriosclerosis, damage to the nerve tissue and for inhibiting the reocclusion of blood vessels after treatment with a balloon catheter, in vascular prosthetics or after the fitting of mechanical devices for holding the blood vessels open (e.g. stents), or other diseases in which cell proliferation or angiogenesis are involved.

By reason of their biological properties the compounds according to the invention may be used on their own or in conjunction with other

pharmacologically active compounds, for example in tumour therapy, in monotherapy or in conjunction with other anti-tumour therapeutic agents, for example in combination with topoisomerase inhibitors (e.g. etoposide), mitosis inhibitors (e.g. vinblastin, taxol), compounds which interact with nucleic acids (e.g. cisplatin, cyclophosphamide, adriamycin), hormone antagonists (e.g. tamoxifen), inhibitors of metabolic processes (e.g. 5-FU etc.), cytokines (e.g. interferons), kinase inhibitors (e.g. EGFR kinase inhibitors such as e.g. Iressa; Gleevec), receptor tyrosine kinase inhibitors with an allosteric effect, antibodies (e.g. herceptin), COX-2 inhibitors, or in conjunction with radiotherapy, etc. These combinations may be administered either simultaneously or sequentially.

The Examples that follow are intended to illustrate the invention more fully:

Example	Name
1.0	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.1	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.2	3-Z-[1-(4-ethoxycarbonyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.3	3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.4	3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.5	3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.6	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-ethylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.7	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propionyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.8	3-Z-[1-(4-(N-(3-dimethylaminopropyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.9	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-butyryl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.10	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-isobutyryl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.11	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

1.12	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-benzoyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.13	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-phenylacetyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.14	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(pyrid-3-yl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.15	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(furan-2-yl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.16	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(2-methoxy-benzoyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.17	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.18	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-isopropylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.19	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-benzylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.20	3-Z-[1-(4-(N-(4-benzyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.21	3-Z-[1-(4-(N-(morpholin-4-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.22	3-Z-[1-(4-(N-(piperidin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.23	3-Z-[1-(4-(N-(benzylmethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.24	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.25	3-Z-[1-(4-((2,6-dimethyl-piperidin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

1.26	3-Z-[1-(4-((N-(2-(2-methoxy-ethoxy)-ethyl)-N-methyl-amino)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.27	3-Z-[1-(4-(triazol-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.28	3-Z-[1-(4-(di-(2-hydroxy-ethyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.29	3-Z-[1-(4-(5-methyl-imidazol-4-yl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.30	3-Z-[1-(4-(N-tert.butoxycarbonyl-ethylamino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.31	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.32	3-Z-[1-(4-(morpholin-4-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.33	3-Z-[1-(4-(N-benzyl-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.34	3-Z-[1-(4-(pyrrolidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.35	3-Z-[1-(4-((4-methyl-piperazin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.36	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-butylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.37	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methoxyacetyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.38	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(3,4-dimethoxy-benzoyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.39	3-Z-[1-(4-(N-(2-hydroxy-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

1.40	3-Z-[1-(4-(N-(2-benzylmethylamino-ethyl)-N-propionyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.41	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(pyrid-4-yl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.42	3-Z-[1-(4-(N-(phthalimido-2-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.43	3-Z-[1-(4-(N-tert.butoxycarbonyl-methylamino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.44	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.45	3-Z-[1-(4-(N-(di-(2-hydroxy-ethyl)-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.46	3-Z-[1-(4-(N-(pyrrolidin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.47	3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.48	3-Z-[1-(4-(N-(imidazol-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.49	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.50	3-Z-[1-(4-(N-(2-(4-tert.butoxycarbonyl-piperazin-1-yl)-ethylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.51	3-Z-[1-(4-(N-(2-benzylmethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.52	3-Z-[1-(4-(4-tert.butoxycarbonyl-piperazin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.53	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-3-methoxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

1.54	3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.55	3-Z-[1-(3-(pyridin-4-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.56	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-pyridin-2-yl-amino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.57	3-Z-[1-(Indazol-6-yl-amino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.58	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-(pyridin-3-yl-amino))-1-phenyl-methylene]-6-chloro-2-indolinone
1.59	trans-3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-amino)-cyclohexylamino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.60	cis-3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-cyclohexylamino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.61	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-3-methyl-pyrrol-3-yl-amino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.62	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl)-3-methyl-pyrrol-3-yl-amino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.63	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.64	3-Z-[1-(4-(2-Diethylamino-ethyl-sulphonyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.65	3-Z-[1-(4-(1-(2-dimethylamino-ethyl)-imidazol-2-yl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.66	3-Z-[1-(4-(piperidin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.67	3-Z-[1-(4-(4-tert.butoxycarbonyl-piperazin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

1.68	3-Z-[1-(4-(N-cyclohexyl-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.69	3-Z-[1-(4-(N-(2-dimethylamino-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.70	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.71	3-Z-[1-(4-(N-(2-(4-methyl-piperazin-1-yl)-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.72	3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.73	3-Z-[1-(4-((4-dimethylamino-piperidin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.74	3-Z-[1-(3-methoxycarbonylmethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.75	3-Z-[1-(3-methoxycarbonyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.76	3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.77	3-Z-[1-(4-(N-tert.butoxycarbonyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.78	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-isopropyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.79	3-Z-[1-(4-(N-(4-tert.butoxycarbonyl-piperazin-1-yl-methylcarbonyl)-N-isopropyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.80	3-Z-[1-(4-(diethylamino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.81	3-Z-[1-(4-(N-propyl-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

1.82	3-Z-[1-(4-(4-methyl-piperazin-1-yl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.83	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.84	3-Z-[1-(4-(N-(4-tert.butoxycarbonyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.85	3-Z-[1-(4-(N-(hydroxy-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.86	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.87	3-Z-[1-(4-(N-(N-(tert.butoxycarbonyl-3-amino-propyl)-N-methyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.88	3-Z-[1-(4-(N-(4-methyl-homopiperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.89	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-3-cyano-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.90	3-Z-[1-(4-(N-(4-ethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.91	3-Z-[1-(4-(N-(1-methyl-piperidin-4-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.92	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.93	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methyl-carbamoylmethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.94	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoylmethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.95	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonylmethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

1.96	3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.97	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-2,3-dimethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.98	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-2,3-dimethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.99	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-aminocarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.100	3-Z-[1-(4-ethoxycarbonylmethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.101	3-Z-[1-(4-(N-(1-methyl-piperidin-4-yl-aminocarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.102	3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.103	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.104	3-Z-[1-(4-(N-(N-(3-dimethylamino-propyl)-aminocarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.105	3-Z-[1-(4-(N-(pyridin-4-yl-methylaminocarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
1.106	3-Z-[1-(4-(N-(1-methyl-piperidin-4-oxy-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
2.0	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-5-nitro-2-indolinone
2.1	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-5-nitro-2-indolinone
2.2	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-5-nitro-2-indolinone

2.3	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-5-nitro-2-indolinone
3.0	3-Z-[1-(4-(N-(dimethyl-carbamoyl-methyl)-N-(2-pyrrolidin-1-yl-ethyl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
3.1	3-Z-[1-(4-(N-(dimethyl-carbamoyl-methyl)-N-(pyrrolidin-1-yl-methyl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
3.2	3-Z-[1-(4-(N-(dimethyl-carbamoyl-methyl)-N-(2-dimethylamino-ethyl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
3.3	3-Z-[1-(4-(N-(dimethyl-carbamoyl-methyl)-N-(dimethylamino-methylcarbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
4.0	3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(3-iodo-phenyl)-methylene]-6-chloro-2-indolinone
4.1	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(3-iodo-phenyl)-methylene]-6-chloro-2-indolinone
4.2	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-chloro-phenyl)-methylene]-6-chloro-2-indolinone
4.3	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-chloro-phenyl)-methylene]-6-chloro-2-indolinone
4.4	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-chloro-phenyl)-methylene]-6-chloro-2-indolinone
4.5	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-chloro-phenyl)-methylene]-6-chloro-2-indolinone
4.6	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-chloro-phenyl)-methylene]-6-chloro-2-indolinone
4.7	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone
4.8	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone

4.9	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone
4.10	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone
4.11	3-Z-[1-(4-(N-(2-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone
4.12	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone
4.13	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone
4.14	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone
4.15	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone
4.16	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone
4.17	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(3-hydroxy-4-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.0	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene]-6-chloro-2-indolinone
5.1	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene]-6-chloro-2-indolinone
5.2	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene]-6-chloro-2-indolinone
5.3	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene]-6-chloro-2-indolinone
5.4	3-Z-[1-(4-(pyrrolidin-1-yl-methyl)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone

5.5	3-Z-[1-(4-(<i>N</i> -(dimethylamino-methylcarbonyl)- <i>N</i> -methyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
5.6	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
5.7	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
5.8	3-Z-[1-(4-(<i>N</i> -(3-dimethylamino-propyl)- <i>N</i> -acetyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
5.9	3-Z-[1-(4-(2-Diethylamino-ethyl-sulphonyl)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
5.10	3-Z-[1-(4-(ethylmethylamino-methyl)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
5.11	3-Z-[1-(4-(<i>N</i> -tert.butoxycarbonyl-methylamino-methyl)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
5.12	3-Z-[1-(4-(<i>N</i> -(dimethylamino-methylcarbonyl)- <i>N</i> -methyl-amino)-anilino)-1-(4-amino-3-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.13	3-Z-[1-(4-(<i>N</i> -(2-dimethylamino-ethyl)- <i>N</i> -methylsulphonyl-amino)-anilino)-1-(4-amino-3-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.14	3-Z-[1-(4-(pyrrolidin-1-yl-methyl)-anilino)-1-(4-amino-3-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.15	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-amino-3-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.16	3-Z-[1-(4-(<i>N</i> -(2-dimethylamino-ethyl)- <i>N</i> -acetyl-amino)-anilino)-1-(4-amino-3-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.17	3-Z-[1-(4-(<i>N</i> -(3-dimethylamino-propyl)- <i>N</i> -acetyl-amino)-anilino)-1-(4-amino-3-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.18	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-nitro-phenyl)-methylene]-6-chloro-2-indolinone

5.19	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.20	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.21	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.22	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-nitro-phenyl)-methylene]-6-chloro-2-indolinone
5.23	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-cyano-phenyl)-methylene]-6-chloro-2-indolinone
6.0	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.1	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.2	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.3	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.4	3-Z-[1-(4-methyl-piperidin-1-yl-amino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.5	3-Z-[1-(4-(N-tert.butoxycarbonyl-ethylamino-methyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.6	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.7	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-benzylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.8	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(n-propylsulphonyl)-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone

6.9	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.10	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.11	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.12	3-Z-[1-(4-(pyrrolidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.13	3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.14	3-Z-[1-(4-((4-methyl-piperazin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.15	3-Z-[1-(4-(N-((2-dimethylamino-ethyl)-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.16	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
6.17	3-Z-[1-(4-(N-tert.butoxycarbonyl-methylamino-methyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
7.0	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.1	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.2	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.3	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.4	3-Z-[1-(4-bromo-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone

7.5	3-Z-[1-(4-(N-tert.butoxycarbonyl-ethylamino-methyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.6	3-Z-[1-(4-(piperidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.7	3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.8	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-carbamoyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.9	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.10	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.11	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.12	3-Z-[1-(4-(pyrrolidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.13	3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.14	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-carbamoyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.15	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.16	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.17	3-Z-[1-(4-(N-((2-dimethylamino-ethyl)-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.18	3-Z-[1-(4-(N-(4-tert.butoxycarbonyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone

7.19	3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.20	3-Z-[1-(3-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.21	3-Z-[1-(4-(N-(morpholin-4-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.22	3-Z-[1-(4-(N-(4-methyl-homopiperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.23	3-Z-[1-(4-(N-(4-ethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.24	3-Z-[1-(3-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.25	3-Z-[1-(3-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
7.26	3-Z-[1-(3-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.27	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methyl-carbamoylmethyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.28	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoylmethyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
7.29	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonylmethyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
8.0	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.1	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.2	3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone

8.3	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.4	3-Z-[1-(4-(N-tert.butoxycarbonyl-ethylamino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.5	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.6	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.7	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.8	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.9	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.10	3-Z-[1-(4-(pyrrolidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.11	3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.12	3-Z-[1-(4-((4-methyl-piperazin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.13	3-Z-[1-(4-(N-((2-dimethylamino-ethyl)-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.14	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.15	3-Z-[1-(4-(piperidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.16	3-Z-[1-(4-(N-tert.butoxycarbonyl-methylamino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone

8.17	3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.18	3-Z-[1-(4-methylsulphonyl-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.19	3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
8.20	3-Z-(1-Anilino-1-phenyl-methylene)-6-fluoro-2-indolinone
8.21	3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
9.0	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-iodo-phenyl)-methylene]-6-fluoro-2-indolinone
9.1	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone
9.2	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone
9.3	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone
9.4	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-(2-acetyl-amino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.5	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-acetyl-amino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.6	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-acetyl-amino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.7	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3,4,5-trifluoro-phenyl)-methylene]-6-fluoro-2-indolinone
9.8	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3,4,5-trifluoro-phenyl)-methylene]-6-fluoro-2-indolinone

9.9	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3,4,5-trifluoro-phenyl)-methylene]-6-fluoro-2-indolinone
9.10	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.11	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-iodo-phenyl)-methylene]-6-fluoro-2-indolinone
9.12	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.13	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.14	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.15	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.16	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-cyanomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.17	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.18	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.19	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-(N-tert.butoxycarbonyl-2-amino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.20	3-Z-[1-(4-(N-acetyl-N-methyl-amino)-anilino)-1-(4-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

9.21	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.22	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.23	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.24	3-Z-[1-(4-(N-tert.butoxycarbonyl-methylamino-methyl)-anilino)-1-(4-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.25	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.26	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(4-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.27	3-Z-[1-(4-methylsulphonyl-anilino)-1-(4-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.28	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.29	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.30	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.31	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.32	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone

9.33	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.34	3-Z-[1-(4-(N-(4-dimethylamino-butylcarbonyl)-N-methyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.35	3-Z-[1-Anilino-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.36	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.37	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.38	3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.39	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.40	3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.41	3-Z-[1-(4-(N-(2-dimethylamino-ethylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.42	3-Z-[1-(4-methylsulphonyl-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.43	3-Z-[1-(4-(N-(3-dimethylamino-propylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.44	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.45	3-Z-[1-Anilino-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone

9.46	3-Z-[1-(4-methylsulphonyl-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.47	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.48	3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.49	3-Z-[1-(4-(N-(3-dimethylamino-propylcarbonyl)-N-methyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.50	3-Z-[1-(4-(N-(2-dimethylamino-ethylcarbonyl)-N-methyl-amino)-anilino)-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
9.51	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.52	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.53	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.54	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.55	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
9.56	3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(3-acetylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone

9.57	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-acetylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
9.58	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-acetylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
9.59	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-acetylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
10.0	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-cyano-2-indolinone
11.0	3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(3-(2-methoxycarbonyl-vinyl)-phenyl)-methylene]-6-chloro-2-indolinone
11.1	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-(2-methoxycarbonyl-vinyl)-phenyl)-methylene]-6-chloro-2-indolinone
11.2	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-(2-carbamoyl-vinyl)-phenyl)-methylene]-6-fluoro-2-indolinone
11.3	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-(2-methoxycarbonyl-vinyl)-phenyl)-methylene]-6-fluoro-2-indolinone
11.4	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-(2-methoxycarbonyl-vinyl)-phenyl)-methylene]-6-fluoro-2-indolinone
12.0	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
12.1	3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
12.2	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-carbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
12.3	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
12.4	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

13.0	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-amino-phenyl)-methylene]-6-chloro-2-indolinone
13.1	3-Z-[1-(4-(N-((2-dimethylamino-ethyl)-carbonyl)-N-methyl-amino)-anilino)-1-(4-amino-phenyl)-methylene]-6-chloro-2-indolinone
13.2	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-amino-phenyl)-methylene]-6-chloro-2-indolinone
13.3	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-amino-phenyl)-methylene]-6-chloro-2-indolinone
13.4	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-amino-phenyl)-methylene]-6-chloro-2-indolinone
14.0	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-aminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
15.0	3-Z-[1-(4-(N-((4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-aminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
16.0	3-Z-[1-(4-(N-((4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone
16.1	3-Z-[1-(4-(N-((2-dimethylamino-ethyl)-carbonyl)-N-methyl-amino)-anilino)-1-(4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone
16.2	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone
16.3	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone
16.4	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone
17.0	3-Z-[1-(4-ethylaminomethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone-trifluoracetat
17.1	3-Z-[1-(4-methylamino-methyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

17.2	3-Z-[1-(4-(N-(2-piperazin-1-yl-ethylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
17.3	3-Z-[1-(4-(piperazin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
17.4	3-Z-[1-(4-(piperazin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
17.5	3-Z-[1-(4-(amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
17.6	3-Z-[1-(4-(N-(piperazin-1-yl-methylcarbonyl)-N-isopropyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
17.7	3-Z-[1-(4-(N-(piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
17.8	3-Z-[1-(4-(N-(N-(3-amino-propyl)-N-methyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
17.9	3-Z-[1-(4-(methylamino-methyl)-anilino)-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
17.10	3-Z-[1-(4-(ethylamino-methyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
17.11	3-Z-[1-(4-(methylamino-methyl)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone
17.12	3-Z-[1-(4-(ethylamino-methyl)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
17.13	3-Z-[1-(4-(piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
17.14	3-Z-[1-(4-(ethylamino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
17.15	3-Z-[1-(4-(methylamino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone

17.16	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-aminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
17.17	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-(2-amino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
17.18	3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-aminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
17.19	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-aminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
17.20	3-Z-[1-(4-(methylamino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
17.21	3-Z-[1-(4-(methylamino-methyl)-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
17.22	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-aminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
18.0	3-Z-[1-(4-(N-aminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
19.0	3-Z-[1-(4-carboxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
19.1	3-Z-[1-(4-carboxymethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
19.2	3-Z-[1-(3-carboxymethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
19.3	3-Z-[1-(3-carboxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
20.0	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
20.1	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.2	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone

20.3	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.4	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.5	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.6	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.7	3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.8	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.9	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.10	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.11	3-Z-[1-(4-(N-tert.butoxycarbonyl-methylamino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.12	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.13	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.14	3-Z-[1-(4-methylsulphonyl-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.15	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.16	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone

20.17	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.18	3-Z-[1-(4-(N-(4-dimethylamino-butylcarbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.19	3-Z-[1-Anilino-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.20	3-Z-[1-(4-methylsulphonyl-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.21	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.22	3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.23	3-Z-[1-Anilino-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.24	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.25	3-Z-[1-(4-(N-(3-dimethylamino-propylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.26	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.27	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.28	3-Z-[1-(4-methylsulphonyl-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.29	3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone

20.30	3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.31	3-Z-[1-(4-(N-(2-dimethylamino-ethylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.32	3-Z-[1-(4-(N-(3-dimethylamino-propylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.33	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.34	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
20.35	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.36	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
20.37	3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.0	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carbamoyl-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
21.1	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
21.2	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.3	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-dimethylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.4	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.5	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

21.6	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-dimethylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.7	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-carbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.8	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-methylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.9	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-carbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.10	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-dimethylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.11	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-(2-(4-methyl-piperazin-1-yl-carbonyl)-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.12	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.13	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-carbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.14	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-dimethylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.15	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-methylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.16	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-methylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.17	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-dimethylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.18	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone

21.19	3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.20	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.21	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.22	3-Z-[1-(4-(N-tert.butoxycarbonyl-methylamino-methyl)-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.23	3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.24	3-Z-[1-(4-methylsulphonyl-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.25	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-(2-methylcarbamoyl-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
21.26	3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(3-methylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
21.27	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-methylcarbamoylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
22.0	3-Z-[1-(4-(4-methyl-piperazin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-1,3-dihydro-indol-2-thion
23.0	3-Z-[1-(4-(N-acetylaminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
23.1	3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-acetyl-amino-phenyl)-methylene]-6-chloro-2-indolinone
23.2	3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-acetyl-amino-phenyl)-methylene]-6-chloro-2-indolinone

23.3	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-acetyl-amino-phenyl)-methylene]-6-chloro-2-indolinone
23.4	3-Z-[1-(4-(piperazin-1-yl-carbonyl)-anilino)-1-(4-acetyl-amino-phenyl)-methylene]-6-chloro-2-indolinone
23.5	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-acetyl-amino-phenyl)-methylene]-6-chloro-2-indolinone
23.6	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-acetylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
23.7	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-acetylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
23.8	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-benzoyl-amino-phenyl)-methylene]-6-chloro-2-indolinone
23.9	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-benzoylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
23.10	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-acetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.11	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-propionylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.12	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-benzoylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.13	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-phenylacetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.14	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-acetyl-amino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.15	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-benzoyl-amino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

23.16	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-propionylamino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.17	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-phenylacetylaminomethyl)-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.18	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-acetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.19	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-propionylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.20	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-phenylacetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.21	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-acetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.22	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-propionylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.23	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-phenylacetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.24	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-cyclopropylcarbonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.25	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-cyclobutylcarbonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.26	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(pyridin-2-yl-carbonylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

23.27	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-cyclohexylcarbonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.28	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(pyridin-3-yl-carbonylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.29	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-isobutyrylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.30	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(3-methylbutyryl-aminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.31	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-cyclohexylmethylcarbonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.32	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-methoxyacetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.33	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-methoxybenzoyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.34	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-tert.butylacetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.35	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-thiophen-carbonylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.36	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-pivaloylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone

23.37	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-furoylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.38	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-acetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.39	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-propionylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.40	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-benzoylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.41	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-phenylacetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.42	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-cyclopropylcarbonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.43	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-cyclobutylcarbonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.44	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(pyridin-2-yl-carbonylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.45	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-cyclohexylcarbonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.46	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(pyridin-3-yl-carbonylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.47	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-isobutyrylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone

23.48	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(3-methylbutyryl-aminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.49	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-cyclohexylmethylcarbonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.50	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-methoxyacetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.51	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-methoxybenzoyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.52	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-tert.butylacetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.53	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-thiophen-carbonylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.54	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-pivaloylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
23.55	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-furoylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
23.56	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(pyridin-4-yl-carbonylaminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
24.0	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-phenylsulphonylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
24.1	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methylsulphonylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
24.2	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-methylsulphonylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone
24.3	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-phenylsulphonylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone

24.4	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-methylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
24.5	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-ethylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
24.6	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-phenylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
24.7	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-ethylsulphonylamino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
24.8	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-methylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
24.9	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-ethylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
24.10	3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-benzylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
24.11	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
24.12	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-ethylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
24.13	3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-benzylsulphonylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone

Abbreviations used:

HOBt = 1-hydroxy-1H-benzotriazole

TBTU = O-benzotriazol-1-yl-N,N,N',N'-tetramethyluronium-tetrafluoroborate

Preparation of the starting compounds:

Example I:

dimethyl 2-(4-fluoro-2-nitrophenyl)-malonate

185 g of potassium-*tert*-butoxide are added to a solution of 188 ml of dimethyl malonate in 970 ml of N-methylpyrrolidone, while cooling with ice, and the mixture is stirred for 2 hours. 150 ml of 2,5-difluoronitrobenzene are added dropwise to the resulting suspension over a period of 30 minutes and then stirred for 6 hours at 85 °C. The mixture is poured onto 4 litres of ice water and 250 ml of concentrated hydrochloric acid and extracted with 2 litres of ethyl acetate. The organic phase is dried with sodium sulphate and evaporated down. The oily residue is stirred twice with water and then taken up in 600 ml ethyl acetate. The solution is dried with sodium sulphate and evaporated to dryness. The crystallised crude product is recrystallised from 600 ml ethyl acetate/hexane = 2:8 and dried.

Yield: 222 g (59 % of theory)

R_f value: 0.40 (silica gel, cyclohexane/ethyl acetate = 5:1)

C₁₁H₁₀FNO₆

Mass spectrum: m/z = 270 [M-H]⁻

The following compounds are prepared analogously to Example I:

(I.1) diethyl 2-(4-bromo-2-nitrophenyl)-malonate

from 2,5-dibromonitrobenzene and diethyl malonate

R_f value: 0.40 (silica gel, petroleum ether/ethyl acetate = 5:1)

C₁₃H₁₄BrNO₆

Mass spectrum: $m/z = 359/361$ $[M]^+$

(1.2) dimethyl 2-(4-cyano-2-nitrophenyl)-malonate

from 4-chloro-3-nitro-benzonitrile and dimethyl malonate

R_f value: 0.50 (silica gel, methylene chloride/methanol = 50:1)

$C_{12}H_{10}N_2O_6$

Mass spectrum: $m/z = 277$ $[M-H]^-$

Example II:

Methyl 4-cyano-2-nitrophenylacetate

14.2 g dimethyl 2-(4-cyano-2-nitrophenyl)-malonate (educt 1.2) are dissolved in 200 ml dimethylsulphoxide and 4.5 g lithium chloride and 1.0 ml of water is added. The solution is stirred for 3.5 hours at 100 °C, then combined with 300 ml ice water and left to stand for 12 hours. The precipitate formed is suction filtered, taken up in methylene chloride and washed with water. The organic phase is dried over sodium sulphate, concentrated in the rotary evaporator and dried.

Yield: 7.7 g (68 % of theory)

R_f value: 0.40 (silica gel, methylene chloride/methanol) = 50:1

$C_{10}H_8N_2O_4$

Mass spectrum: $m/z = 219$ $[M-H]^-$

Example III:

4-fluoro-2-nitrophenylacetic acid

50.0 g of dimethyl 2-(4-fluoro-2-nitrophenyl)-malonate (educt I) are stirred in 400 ml of 6 molar hydrochloric acid for 20 hours at 100 °C, then combined with 400 ml of water and cooled to 0 °C. The precipitate formed is suction filtered, washed with water and 100 ml petroleum ether and dried.

Yield: 34.5 g (94 % of theory)

R_f value: 0.30 (silica gel, cyclohexane/ethyl acetate) = 5:2

$C_8H_6FNO_4$

Mass spectrum: $m/z = 154$ $[M-COO-H]^-$

Example IV:

6-fluoro-2-indolinone

119 g 4-fluoro-2-nitrophenylacetic acid (educt III) are hydrogenated in 600 ml acetic acid with the addition of 20 g palladium on activated charcoal (10%) under 50 psi of hydrogen pressure. The catalyst is suction filtered and the solvent is distilled off. The crude product is extracted with 500 ml petroleum ether, suction filtered, washed with water and dried.

Yield: 82.5 g (91 % of theory)

R_f value: 0.30 (silica gel, petroleum ether/ethyl acetate = 1:1)

C₈H₆FNO

Mass spectrum: m/z = 150 [M-H]⁻

The following compounds are prepared analogously to Example IV:

(IV.1) 6-bromo-2-indolinone

from diethyl 2-(4-bromo-2-nitrophenyl)-malonate (educt I.1) with Raney nickel as hydrogenation catalyst

R_f value: 0.45 (silica gel, petroleum ether/ethyl acetate = 1:1)

C₈H₆BrNO

Mass spectrum: m/z = 210/212 [M-H]⁻

(IV.2) 6-cyano-2-indolinone

from methyl 4-cyano-2-nitrophenylacetate (educt II) with palladium/calcium carbonate as hydrogenation catalyst

R_f value: 0.45 (silica gel, methylene chloride/methanol = 9:1)

C₉H₆N₂O

Mass spectrum: m/z = 157 [M-H]⁻

Example V:

1-acetyl-6-fluoro-2-indolinone

82.5 g 6-fluoro-2-indolinone (educt IV) are stirred in 180 ml acetic anhydride for 3 hours at 130 °C. After cooling to ambient temperature the precipitate is suction filtered, washed with 100 ml of petroleum ether and dried.

Yield: 64.8 g (61 % of theory)

R_f value: 0.75 (silica gel, petroleum ether/ethyl acetate = 1:1)

C₁₀H₈FNO₂

Mass spectrum: m/z = 192 [M-H]⁻

The following compounds are prepared analogously to Example V:

(V.1) 1-acetyl-6-chloro-2-indolinone

from 6-chloro-2-indolinone and acetic anhydride

R_f value: 0.55 (silica gel, petroleum ether/ethyl acetate = 2:3)

C₁₁H₁₀ClNO₂

Mass spectrum: m/z = 208/210 [M-H]⁻

(V.2) 1-acetyl-6-bromo-2-indolinone

from 6-bromo-2-indolinone (educt IV.1) and acetic anhydride

R_f value: 0.60 (silica gel, petroleum ether/ethyl acetate = 2:1)

C₁₀H₈BrNO₂

Mass spectrum: m/z = 253/255 [M]⁺

(V.3) 1-acetyl-6-cyano-2-indolinone

from 6-cyano-2-indolinone (educt IV.2) and acetic anhydride

R_f value: 0.60 (silica gel, methylene chloride/methanol = 50:1)

C₁₁H₈N₂O₂

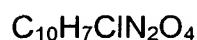
Mass spectrum: m/z = 199 [M-H]⁻

Example VI:

1-acetyl-5-nitro-6-chloro-2-indolinone

2.75 g 1-acetyl-6-chloro-2-indolinone (educt V.1) are placed in 40 ml concentrated sulphuric acid and at -10°C 1.05 g of ammonium nitrate are added. The mixture is stirred for 1.5 hours at ambient temperature. After this time the mixture is poured onto ice water and stirred for a further 20 minutes. The solution is neutralised with concentrated ammonia, the precipitate formed is suction filtered and washed with a little ethanol and ether.

Yield: 2.80 g (84 % of theory)



Mass spectrum: $m/z = 254$ $[\text{M}]^+$

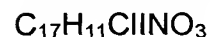
Example VII:

1-acetyl-3-[1-hydroxy-1-(3-iodo-phenyl)-methylene]-6-chloro-2-indolinone

10.5 g 1-acetyl-6-chloro-2-indolinone (educt V.1), 13.6 g 3-iodobenzoic acid and 17.7 g TBTU are placed in 100 ml of dimethylformamide, 35 ml triethylamine are added and the mixture is stirred for 12 hours at ambient temperature. After this time the solvent is eliminated, the residue is combined with water, suction filtered and washed with a little water, methanol and ether and dried in vacuo at 100°C .

Yield: 12.9 g (59 % of theory)

R_f value: 0.80 (silica gel, methylene chloride/methanol = 9:1)



Mass spectrum: $m/z = 438/440$ $[\text{M}-\text{H}]^-$

The following compounds are prepared analogously to Example VII:

(VII.1) 1-acetyl-3-[1-hydroxy-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and methyl (4-carboxyphenyl)-acetate (prepared according to Tetrahedron **1997**, 53, 7335-7340)

(VII.2) 1-acetyl-3-[1-hydroxy-1-(4-chloro-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 4-chloro-benzoic acid

(VII.3) 1-acetyl-3-[1-hydroxy-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 4-benzyloxy-benzoic acid

(VII.4) 1-acetyl-3-[1-hydroxy-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 3,4-dimethoxy-benzoic acid

(VII.5) 1-acetyl-3-[1-hydroxy-1-(3,4-dimethoxy-phenyl)-methylene]-6-cyano-2-indolinone

from 1-acetyl-6-cyano-2-indolinone (educt V.3) and 3,4-dimethoxy-benzoic acid

(VII.6) 1-acetyl-3-[1-hydroxy-1-(3-nitro-4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 3-nitro-4-hydroxy-benzoic acid

(VII.7) 1-acetyl-3-[1-hydroxy-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 3-fluoro-benzoic acid

(VII.8) 1-acetyl-3-[1-hydroxy-1-(4-(2-acetylamino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 4-(2-acetylamino-ethyl)-benzoic acid (prepared according to J. Am. Chem. Soc. **1943**, 65, 2377)

(VII.9) 1-acetyl-3-[1-hydroxy-1-(3,4,5-trifluoro-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 3,4,5-trifluoro-benzoic acid

(VII.10) 1-acetyl-3-[1-hydroxy-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and (3-carboxyphenyl)-acetate methyl (prepared analogously to Tetrahedron **1997**, 53, 7335-7340)

(VII.11) 1-acetyl-3-[1-hydroxy-1-(3-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 3-(N-tert.butoxycarbonyl-aminomethyl)-benzoic acid (prepared according to Tetrahedron **1997**, 53, 7335-7340)

(VII.12) 1-acetyl-3-[1-hydroxy-1-(3-cyanomethyl-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and (3-carboxy-phenyl)-acetonitrile (prepared according to J. Prakt. Chem. **1998**, 340, 367-374)

(VII.13) 1-acetyl-3-[1-hydroxy-1-(4-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 4-(N-tert.butoxycarbonyl-aminomethyl)-benzoic acid (prepared according to Bioorg. Med. Chem. Lett **2000**, 10, 553-557)

(VII.14) 1-acetyl-3-[1-hydroxy-1-(4-iodo-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 4-iodo-benzoic acid

(VII.15) 1-acetyl-3-[1-hydroxy-1-(4-iodo-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 4-iodo-benzoic acid

(VII.16) 1-acetyl-3-[1-hydroxy-1-(3-iodo-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 3-iodo-benzoic acid

(VII.17) 1-acetyl-3-[1-hydroxy-1-(4-(2-methoxycarbonylethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 4-(2-methoxycarbonylethyl)-benzoic acid (prepared analogously to Tetrahedron **1997**, 53, 7335-7340)

(VII.18) 1-acetyl-3-[1-hydroxy-1-(3-(2-methoxycarbonylethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 3-(2-methoxycarbonylethyl)-benzoic acid (prepared analogously to Tetrahedron **1997**, 53, 7335-7340)

(VII.19) 1-acetyl-3-[1-hydroxy-1-(3-(N-tert.butoxycarbonyl-2-aminoethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 3-(N-tert.butoxycarbonyl-2-aminoethyl)-benzoic acid (prepared analogously to Bioorg. Med. Chem. Lett **2000**, 10, 553-557)

(VII.20) 1-acetyl-3-[1-hydroxy-1-(4-(N-tert.butoxycarbonyl-2-aminoethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-6-fluoro-2-indolinone (educt V) and 4-(N-tert.butoxycarbonyl-2-aminoethyl)-benzoic acid (prepared analogously to Bioorg. Med. Chem. Lett **2000**, 10, 553-557)

(VII.21) 1-acetyl-3-[1-hydroxy-1-(4-acetylamino-3-nitro-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 3-acetylamino-4-nitro-benzoic acid

(VII.22) 1-acetyl-3-[1-hydroxy-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 4-(imidazol-1-yl-methyl)-benzoic acid (prepared according to J. Med. Chem. **1987**, 30, 1342-1347)

(VII.23) 1-acetyl-3-[1-hydroxy-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone
 from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 4-(2-oxo-pyrrolidin-1-yl)-benzoic acid (prepared according to J. Med. Chem. **1999**, 42, 2332-2343)

(VII.24) 1-acetyl-3-[1-hydroxy-1-(4-nitro-phenyl)-methylene]-6-chloro-2-indolinone
 from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 4-nitro-benzoic acid

(VII.25) 1-acetyl-3-[1-hydroxy-1-(4-cyano-phenyl)-methylene]-6-chloro-2-indolinone
 from 1-acetyl-6-chloro-2-indolinone (educt V.1) and 4-cyano-benzoic acid

(VII.26) 1-acetyl-3-[1-hydroxy-1-(3-acetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
 from 1-acetyl-6-fluoro-2-indolinone (educt V) and 3-acetylaminomethyl-benzoic acid (prepared according to J. Med. Chem. **1997**, 40, 4030-4052)

Example VIII:

1-acetyl-3-[1-methoxy-1-(3-iodo-phenyl)-methylene]-6-chloro-2-indolinone

2.36 g of trimethyloxonium tetrafluoroborate are added batchwise to a solution of 3.52 g of 1-acetyl-3-[1-hydroxy-1-(3-iodo-phenyl)-methylene]-6-chloro-2-indolinone (educt VII) and 2.72 ml of ethyldiisopropylamine in 80 ml dichloromethane and the mixture is stirred for one hour at ambient temperature. Then another 1.4 ml ethyldiisopropylamine and 1.2 g trimethyloxonium tetrafluoroborate are added and the mixture is stirred for a further two hours at ambient temperature. Then it is extracted with water, the organic phase is dried over magnesium sulphate and evaporated to dryness. The residue is recrystallised from ether and dried in vacuo at 80 °C.

Yield: 2.40 g (66 % of theory)

R_f value: 0.60 (silica gel, petroleum ether/dichloromethane/ethyl acetate = 5:4:1)

C₁₈H₁₃ClINO₃

Mass spectrum: $m/z = 438/440$ $[M-H]^-$

M.p. 185 – 187 °C

The following compounds are prepared analogously to Example VIII:

(VIII.1) 1-acetyl-3-[1-methoxy-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone
from 1-acetyl-3-[1-hydroxy-1-(4-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.1)

(VIII.2) 1-acetyl-3-[1-methoxy-1-(4-chloro-phenyl)-methylene]-6-chloro-2-indolinone
from 1-acetyl-3-[1-hydroxy-1-(4-chloro-phenyl)-methylene]-6-chloro-2-indolinone (educt VII.2)

(VIII.3) 1-acetyl-3-[1-methoxy-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone
from 1-acetyl-3-[1-hydroxy-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone (educt VII.3)

(VIII.4) 1-acetyl-3-[1-methoxy-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone
from 1-acetyl-3-[1-hydroxy-1-(3,4-dimethoxy-phenyl)-methylene]-6-chloro-2-indolinone (educt VII.4)

(VIII.5) 1-acetyl-3-[1-methoxy-1-(3,4-dimethoxy-phenyl)-methylene]-6-cyano-2-indolinone
from 1-acetyl-3-[1-hydroxy-1-(3,4-dimethoxy-phenyl)-methylene]-6-cyano-2-indolinone (educt VII.5)

(VIII.6) 1-acetyl-3-[1-methoxy-1-(3-nitro-4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone
from 1-acetyl-3-[1-hydroxy-1-(3-nitro-4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone (educt VII.6)

(VIII.7) 1-acetyl-3-[1-methoxy-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(3-fluoro-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.7)

(VIII.8) 1-acetyl-3-[1-methoxy-1-(4-(2-acetylamino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(4-(2-acetylamino-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.8)

(VIII.9) 1-acetyl-3-[1-methoxy-1-(3,4,5-trifluoro-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(3,4,5-trifluoro-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.9)

(VIII.10) 1-acetyl-3-[1-methoxy-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(3-methoxycarbonylmethyl-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.10)

(VIII.11) 1-acetyl-3-[1-methoxy-1-(3-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(3-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.11)

(VIII.12) 1-acetyl-3-[1-methoxy-1-(3-cyanomethyl-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(3-cyanomethyl-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.12)

(VIII.13) 1-acetyl-3-[1-methoxy-1-(4-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(4-(N-tert.butoxycarbonyl-aminomethyl)-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.13)

(VIII.14) 1-acetyl-3-[1-methoxy-1-(4-iodo-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(4-iodo-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.14)

(VIII.15) 1-acetyl-3-[1-methoxy-1-(4-iodo-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(4-iodo-phenyl)-methylene]-6-chloro-2-indolinone (educt VII.15)

(VIII.16) 1-acetyl-3-[1-methoxy-1-(3-iodo-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(3-iodo-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.16)

(VIII.17) 1-acetyl-3-[1-methoxy-1-(3-(2-methoxycarbonylethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(3-(2-methoxycarbonylethyl)-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.17)

(VIII.18) 1-acetyl-3-[1-methoxy-1-(4-(2-methoxycarbonylethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(4-(2-methoxycarbonylethyl)-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.18)

(VIII.19) 1-acetyl-3-[1-methoxy-1-(4-(N-tert.butoxycarbonyl-2-aminoethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(4-(N-tert.butoxycarbonyl-2-aminoethyl)-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.20)

(VIII.20) 1-acetyl-3-[1-methoxy-1-(3-(N-tert.butoxycarbonyl-2-aminoethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
 from 1-acetyl-3-[1-hydroxy-1-(3-(N-tert.butoxycarbonyl-2-aminoethyl)-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.19)

(VIII.21) 1-acetyl-3-[1-methoxy-1-(3-acetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone
 from 1-acetyl-3-[1-hydroxy-1-(3-acetylaminomethyl-phenyl)-methylene]-6-fluoro-2-indolinone (educt VII.26)

Example IX:

1-acetyl-3-[1-ethoxy-1-phenyl-methylene]-6-chloro-2-indolinone

A solution of 41.9 g 1-acetyl-6-chloro-2-indolinone (educt V.1) and 136 ml triethyl orthobenzoate in 150 ml acetic anhydride is stirred for six hours at 120°C. After cooling the mixture is evaporated down by half, the crystals precipitated are suction filtered and dried at 100 °C in vacuo.

Yield: 38.0 g (56 % of theory)

R_f value: 0.60 (silica gel, petroleum ether/dichloromethane/ethyl acetate = 5:4:1)

C₁₉H₁₆ClNO₃

Mass spectrum: m/z = 342/344 [M+H]⁺

M.p. 185 – 187 °C

The following compounds are prepared analogously to Example IX:

(IX.1) 1-acetyl-3-[1-ethoxy-1-phenyl-methylene]-6-bromo-2-indolinone
 from 1-acetyl-6-bromo-2-indolinone (educt V.2)

(IX.2) 1-acetyl-3-[1-ethoxy-1-phenyl-methylene]-6-cyano-2-indolinone
 from 1-acetyl-6-cyano-2-indolinone (educt V.3)

(IX.3) 1-acetyl-3-[1-ethoxy-1-phenyl-methylene]-6-fluoro-2-indolinone
 from 1-acetyl-6-fluoro-2-indolinone (educt V)

(IX.4) 1-acetyl-3-[1-ethoxy-1-phenyl-methylene]-5-nitro-6-chloro-2-indolinone
from 1-acetyl-5-nitro-6-chloro-2-indolinone (educt VI)

Example X:

1-acetyl-3-[1-chloro-1-(4-acetylamino-3-nitro-phenyl)methylene]-2-indolinone

A suspension of 11.4 g of 1-acetyl-3-(1-hydroxy-1-(4-acetylamino-3-nitro-phenyl)methylene)-2-indolinone (educt VII.21) and 9.37 g of phosphorus pentachloride in 200 ml dioxane is stirred for 4 hours at 100 °C. After the addition of a further 2.0 g of phosphorus pentachloride the mixture is stirred for a further 3 hours at 100 °C. Then the solvent is distilled off, the residue is stirred with 100 ml of ethyl acetate, suction filtered, washed with ethyl acetate and dried at 60 °C.

Yield: 6.40 g (53 % of theory)

R_f value: 0.70 (silica gel, dichloromethane/ethyl acetate = 9:1)

C₁₉H₁₄ClN₃O₅

Mass spectrum: m/z = 398/400 [M-H]⁻

The following compounds are prepared analogously to Example X:

(X.1) 1-acetyl-3-[1-chloro-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-3-[1-hydroxy-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene]-6-chloro-2-indolinone (educt VII.22)

(X.2) 1-acetyl-3-[1-chloro-1-(4-(2-oxo-pyrrolidin-1-yl)-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-3-[1-chloro-1-(4-(2-oxo-pyrrolidin-1-yl)-methyl)-phenyl)-methylene]-6-chloro-2-indolinone (educt VII.23)

(X.3) 1-acetyl-3-[1-chloro-1-(4-nitro-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-3-[1-chloro-1-(4-nitro-methyl)-phenyl)-methylene]-6-chloro-2-indolinone (educt VII.24)

(X.4) 1-acetyl-3-[1-chloro-1-(4-cyano-phenyl)-methylene]-6-chloro-2-indolinone

from 1-acetyl-3-[1-chloro-1-(4-cyano-methyl)-phenyl]-methylene]-6-chloro-2-indolinone (educt VII.25)

Example XI:

N-(2-dimethylamino-ethyl)-4-nitro-benzamide

1.25 ml N,N-dimethylaminoethylamine are dissolved with 3 ml triethylamine in 20 ml methylene chloride and cooled to 0°C. Then 2 g of 4-nitrobenzoic acid chloride are added batchwise and the mixture is stirred for 5 minutes in the cold and for 20 minutes at ambient temperature. Finally, the precipitate is removed by suction filtering and the organic phase is washed with water, dried over sodium sulphate and concentrated in the rotary evaporator.

Yield: 1.8 g (70 % of theory)

R_f value: 0.78 (silica gel, methylene chloride/methanol = 9:1)

C₁₁H₁₅N₃O₃

Mass spectrum: m/z = 238 [M+H]⁺

The following compounds are prepared analogously to Example XI:

(XI.1) N-(2-dimethylamino-ethyl)-N-methyl-4-nitro-benzamide

(XI.2) N-(3-dimethylamino-propyl)-4-nitro-benzamide

(XI.3) N-(3-dimethylamino-propyl)-N-methyl-4-nitro-benzamide

(XI.4) N-(2-dimethylamino-ethyl)-N-ethyl-4-nitro-benzamide.

(XI.5) N-(2-(tert-butyloxycarbonyl-methylamino-ethyl)-N-methyl-4-nitro-benzamide

- (XI.6) N,N-bis-(2-diethylamino-ethyl)-4-nitro-benzamide
- (XI.7) N-(2-tert-butyloxycarbonyl-amino-ethyl)-4-nitro-benzamide
- (XI.8) N-(2-dimethylamino-ethyl)-3-nitro-benzamide
- (XI.9) N-(2-dimethylamino-ethyl)-N-methyl-3-nitro-benzamide
- (XI.10) N-(3-dimethylamino-propyl)-3-nitro-benzamide
- (XI.11) N-(3-dimethylamino-propyl)-N-methyl-3-nitro-benzamide
- (XI.12) 2-N-(dimethylamino-methyl)-carbamoyl-5-nitro-furan
- (XI.13) 4-(4-methyl-piperazin-1-yl-carbonyl)-nitrobenzene
- (XI.14) 4-(piperidin-1-yl-carbonyl)-nitrobenzene
- (XI.15) N-cyclohexyl-N-methyl-4-nitro-benzamide
- (XI.16) N-isopropyl-4-nitro-benzamide
- (XI.17) 4-(2,3,4,5-tetrahydro-1(H)-benzo[d]azepin-3-yl-carbonyl)-nitrobenzene
- (XI.18) 4-(4-hydroxy-piperidin-1-yl-carbonyl)-nitrobenzene
- (XI.19) 4-(4-tert-butyloxycarbonyl-piperazin-1-yl-carbonyl)-nitrobenzene
- (XI.20) 4-(4-tert-butyloxycarbonyl-[1,4]diazepan-1-yl-carbonyl)-nitrobenzene
- (XI.21) 4-(N-(3-dimethylamino-propyl)-N-methyl-carbamoylmethyl)-nitrobenzene

(XI.22) 4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoylmethyl)-nitrobenzene

(XI.23) 4-[(4-methyl-piperazin-1-yl)-carbonylmethyl]-nitrobenzene

Example XII:

2-amino-5-(4-methyl-piperazin-1-yl-carbonyl)-pyridine

3.00 g 6-amino-nicotinic acid are dissolved in 30 ml of dimethylformamide and 4.05 g of N,N-carbonyldiimidazole are added. The mixture is briefly heated to 70°C and then stirred for a further hour at ambient temperature. After this time 4.85 ml of N-methylpiperazine are added and the mixture is stirred for 12 hours at ambient temperature. The solvent is eliminated and the residue is purified through a silica gel column with methylene chloride/ethanol/ammonia 7:1:0.1 as eluant.

Yield: 4.1 g (86 % of theory)

R_f value: 0.60 (silica gel, methylene chloride/ethanol/ammonia = 5:1:0.1)

C₁₁H₁₆N₄O

Mass spectrum: m/z = 221 [M+H]⁺

The following compounds are prepared analogously to Example XII:

(XII.1) 4-nitro-1-methyl-2-[(2-dimethylamino-ethyl)-N-methyl-carbamoyl]-pyrrole

(XII.2) 4-nitro-1-methyl-2-[(4-methyl-piperazin-1-yl)-carbonyl]-pyrrole

Example XIII:

4-(dimethylamino-ethoxy)-nitrobenzene

5.2 g p-nitrophenol are dissolved in 200 ml acetone and 7.2 g 2-chloro-N,N-dimethylethylamin-hydrochloride and 11 g potassium carbonate are added. The mixture is stirred for 12 hours at reflux temperature. After cooling the salts are filtered off, the filtrate is evaporated down and the

residue is taken up in methylene chloride. The organic phase is washed with water, dried over sodium sulphate and finally the solvent is eliminated.

Yield: 4.1 g (53% of theory)

R_f value: 0.45 (silica gel, methylene chloride/methanol = 9:1)

C₁₀H₁₄N₂O₃

Mass spectrum: m/z = 211 [M+H]⁺

Example XIV:

4-{N-[(4-methyl-piperazin-1-yl)-aminocarbonyl]-N-methyl-amino}-nitrobenzene

11.6 ml of phosgene solution in toluene (20%) are dissolved in 20 ml of tetrahydrofuran and 3.0 g of N-methyl-4-nitroaniline and 2.8 ml triethylamine in 30 ml of tetrahydrofuran are added dropwise at 0 °C. The mixture is stirred for 0.5 hours at 0 °C and for a further hour at ambient temperature. Then it is cooled again to 0°C, 5.4 ml of 1-amino-4-methyl-piperazine in 10 ml of tetrahydrofuran are added and the mixture is stirred for 2 hours at ambient temperature. After this time the solvent is eliminated, the residue is taken up in ethyl acetate and extracted with water. The organic phase is dried over sodium sulphate and finally the solvent is removed. The residue is purified through an aluminium oxide column (activity 2-3) with methylene chloride/ethanol 40:1 as eluant.

Yield: 3.3 g (56% of theory)

R_f value: 0.30 (aluminium oxide, methylene chloride/ethanol = 40:1)

M.p. 170-172 °C

C₁₃H₁₉N₅O₃

Mass spectrum: m/z = 294 [M+H]⁺

The following compounds are prepared analogously to Example XIV:

(XIV.1) 4-{N-[(1-methyl-piperidin-4-yl)-aminocarbonyl]-N-methyl-amino}-nitrobenzene

(XIV.2) 4-{N-[(4-methyl-piperazin-1-yl)-carbonyl]-N-methyl-amino}-nitrobenzene

(XIV.3) 4-{N-[N-(3-dimethylamino-propyl)-aminocarbonyl]-N-methyl-amino}-nitrobenzene

(XIV.4) 4-[N-(pyridin-4-yl-methylaminocarbonyl)-N-methyl-amino]-nitrobenzene

(XIV.5) 4-{N-[(1-methyl-piperidin-4-oxy)-carbonyl]-N-methyl-amino}-nitrobenzene

Example XV:

The syntheses of the following compounds are already described in International Application WO 01/27081:

(XV.1) 4-[(2,6-dimethyl-piperidin-1-yl)-methyl]-aniline

(XV.2) N-(2-dimethylamino-ethyl)-N-methylsulphonyl-p-phenylenediamine

(XV.3) 3-(dimethylaminomethyl)-aniline

(XV.4) 4-(dimethylaminomethyl)-aniline

(XV.5) 4-(2-dimethylamino-ethyl)-aniline

(XV.6) 4-[N-(2-dimethylamino-ethyl)-N-acetyl-amino]-aniline

(XV.7) 4-[N-(3-dimethylamino-propyl)-N-acetyl-amino]-aniline

(XV.8) 4-[N-(2-dimethylamino-ethyl)-N-benzoyl-amino]-aniline

(XV.9) 4-[N-(2-dimethylamino-ethyl)-N-propionyl-amino]-aniline

(XV.10) 4-[N-(2-dimethylamino-ethyl)-N-butyryl-amino]-aniline

(XV.11) 4-[N-(2-dimethylamino-ethyl)-N-isobutyryl-amino]-aniline

(XV.12) 4-(N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.13) 4-(N-ethyl-N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.14) 4-[N-(4-chlorophenyl-methyl)-N-tert.butoxycarbonyl-aminomethyl]-aniline

(XV.15) 4-(N-cyclohexyl-N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.16) 4-(N-isopropyl-N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.17) 4-(N-propyl-N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.18) 4-(N-methyl-N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.19) 4-(N-butyl-N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.20) 4-(N-methoxycarbonyl-methyl-N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.21) 4-(N-benzyl-N-tert.butoxycarbonyl-aminomethyl)-aniline

(XV.22) 4-(pyrrolidin-1-yl-methyl)-aniline

(XV.23) 4-(morpholin-4-yl-methyl)-aniline

(XV.24) 4-(hexamethyleneiminomethyl)-aniline

(XV.25) 4-(4-hydroxy-piperidin-1-yl-methyl)-aniline

(XV.26) 4-(4-methoxy-piperidin-1-yl-methyl)-aniline

(XV.27) 4-(4-methyl-piperidin-1-yl-methyl)-aniline

(XV.28) 4-(4-ethyl-piperidin-1-yl-methyl)-aniline

(XV.29) 4-(4-isopropyl-piperidin-1-yl-methyl)-aniline

(XV.30) 4-(4-phenyl-piperidin-1-yl-methyl)-aniline

(XV.31) 4-(4-benzyl-piperidin-1-yl-methyl)-aniline

(XV.32) 4-(4-ethoxycarbonyl-piperidin-1-yl-methyl)-aniline

(XV.33) 4-(N,N-dipropyl-aminomethyl)-aniline

(XV.34) 4-(4-tert.butoxycarbonyl-piperazin-1-yl-methyl)-aniline

(XV.35) 4-(2-morpholin-4-yl-ethyl)-aniline

(XV.36) 4-(2-pyrrolidin-1-yl-ethyl)-aniline

(XV.37) 4-(2-piperidin-1-yl-ethyl)-aniline

(XV.38) 4-(N-propyl-N-benzyl-aminomethyl)-aniline

(XV.39) 4-[N-(n-hexyl)-N-methyl-aminomethyl]-aniline

(XV.40) 4-[N-methyl-N-(4-chlorbenzyl)-aminomethyl]-aniline

(XV.41) 4-[N-methyl-N-(4-brombenzyl)-aminomethyl]-aniline

(XV.42) 4-[N-methyl-N-(4-methylbenzyl)-aminomethyl]-aniline

(XV.43) 4-[N-methyl-N-(4-fluorbenzyl)-aminomethyl]-aniline

- (XV.44) 4-[N-methyl-N-(3-chlorobenzyl)-aminomethyl]-aniline
- (XV.45) 4-[N-methyl-N-(3,4-dimethoxybenzyl)-aminomethyl]-aniline
- (XV.46) 4-[N-methyl-N-(4-methoxybenzyl)-aminomethyl]-aniline
- (XV.47) 4-(N-2,2,2-trifluorethyl-N-benzyl-aminomethyl)-aniline
- (XV.48) 4-[N-2,2,2-trifluorethyl-N-(4-chlorobenzyl)-aminomethyl]-aniline
- (XV.49) 4-(thiomorpholin-4-yl-methyl)-aniline
- (XV.50) 4-(1-oxo-thiomorpholin-4-yl-methyl)-aniline
- (XV.51) 4-(1,1-dioxo-thiomorpholin-4-yl-methyl)-aniline
- (XV.52) 4-(azetidin-1-yl-methyl)-aniline
- (XV.53) 4-(3,4-dihydropyrrolidin-1-yl-methyl)-aniline
- (XV.54) 4-(3,4-dihydropiperidin-1-yl-methyl)-aniline
- (XV.55) 4-(2-methoxycarbonyl-pyrrolidin-1-yl-methyl)-aniline
- (XV.56) 4-(3,5-dimethyl-piperidin-1-yl-methyl)-aniline
- (XV.57) 4-(4-phenyl-piperazin-1-yl-methyl)-aniline
- (XV.58) 4-(4-phenyl-4-hydroxy-piperidin-1-yl-methyl)-aniline
- (XV.59) 4-[N-(3,4,5-trimethoxy-benzyl)-N-methyl-aminomethyl]-aniline
- (XV.60) 4-[N-(3,4-dimethoxy-benzyl)-N-ethyl-aminomethyl]-aniline

(XV.61) 4-(N-benzyl-N-ethyl-aminomethyl)-aniline

(XV.62) 4-[N-(2,6-dichlorobenzyl)-N-methyl-aminomethyl]-aniline

(XV.63) 4-[N-(4-trifluoromethylbenzyl)-N-methyl-aminomethyl]-aniline

(XV.64) 4-(N-benzyl-N-isopropyl-aminomethyl)-aniline

(XV.65) 4-(N-benzyl-N-tert.butyl-aminomethyl)-aniline

(XV.66) 4-(diethylamino-methyl)-aniline

(XV.67) 4-(2-diethylamino-ethyl)-aniline

(XV.68) 4-(N,N-diisopropyl-aminomethyl)-aniline

(XV.69) 4-(N,N-diisobutyl-aminomethyl)-aniline

(XV.70) 4-(2,3,4,5-tetrahydro-benzo(d)azepin-3-yl-methyl)-aniline

(XV.71) 4-(2,3-dihydro-isoindol-2-yl-methyl)-aniline

(XV.72) 4-(6,7-dimethoxy-1,2,3,4-tetrahydro-isoquinolin-2-yl-methyl)-aniline

(XV.73) 4-(1,2,3,4-tetrahydro-isoquinolin-2-yl-methyl)-aniline

(XV.74) 4-[N-(2-hydroxy-ethyl)-N-benzyl-aminomethyl]-aniline

(XV.75) 4-[N-(1-ethyl-pentyl)-N-(pyridin-2-yl-methyl)-aminomethyl]-aniline

(XV.76) 4-(piperidin-1-yl-methyl)-3-nitro-aniline

(XV.77) 4-(piperidin-1-yl-methyl)-3-amino-aniline

(XV.78) 4-(N-benzyl-N-methyl-aminomethyl)-aniline

(XV.79) 4-(N-ethyl-N-methyl-aminomethyl)-aniline

(XV.80) 4-(N-phenethyl-N-methyl-aminomethyl)-aniline

(XV.81) 4-[N-(3,4-dihydroxy-phenethyl)-N-methyl-aminomethyl]-aniline

(XV.82) 4-[N-(3,4,5-trimethoxy-phenethyl)-N-methyl-aminomethyl]-aniline

(XV.83) 4-[N-(3,4-dimethoxy-phenethyl)-N-methyl-aminomethyl]-aniline

(XV.84) 4-[N-(3,4-dimethoxy-benzyl)-N-methyl-aminomethyl]-aniline

(XV.85) 4-[N-(4-chloro-benzyl)-N-methyl-aminomethyl]-aniline

(XV.86) 4-[N-(4-bromo-benzyl)-N-methyl-aminomethyl]-aniline

(XV.87) 4-[N-(4-fluoro-benzyl)-N-methyl-aminomethyl]-aniline

(XV.88) 4-[N-(4-methyl-benzyl)-N-methyl-aminomethyl]-aniline

(XV.89) 4-[N-(4-nitro-phenethyl)-N-methyl-aminomethyl]-aniline

(XV.90) 4-(N-phenethyl-N-benzyl-aminomethyl)-aniline

(XV.91) 4-(N-phenethyl-N-cyclohexyl-aminomethyl)-aniline

(XV.92) 4-[N-(2-(pyridin-2-yl)-ethyl)-N-methyl-aminomethyl]-aniline

(XV.93) 4-[N-(2-(pyridin-4-yl)-ethyl)-N-methyl-aminomethyl]-aniline

(XV.94) 4-[N-(pyridin-4-yl-methyl)-N-methyl-aminomethyl]-aniline

(XV.95) 4-(N,N-dibenzylaminomethyl)-aniline

(XV.96) 4-[N-(4-nitro-benzyl)-N-propyl-aminomethyl]-aniline

(XV.97) 4-[N-benzyl-N-(3-cyano-propyl)-aminomethyl]-aniline

(XV.98) 4-(N-benzyl-N-allyl-aminomethyl)-aniline

(XV.99) 4-[N-benzyl-N-(2,2,2-trifluorethyl)-aminomethyl]-aniline

(XV.100) 4-[(benzo(1,3)dioxol-5-yl-methyl)-methyl-aminomethyl]-aniline

(XV.101) 4-(7-chloro-2,3,4,5-tetrahydro-benzo(d)azepin-3-yl-methyl)-aniline

(XV.102) 4-(7,8-dichloro-2,3,4,5-tetrahydro-benzo(d)azepin-3-yl-methyl)-
aniline

(XV.103) 4-(7-methoxy-2,3,4,5-tetrahydro-benzo(d)azepin-3-yl-methyl)-aniline

(XV.104) 4-(7-methyl-2,3,4,5-tetrahydro-benzo(d)azepin-3-yl-methyl)-aniline

(XV.105) 4-(7,8-dimethoxy-2,3,4,5-tetrahydro-benzo(d)azepin-3-yl-methyl)-
aniline

(XV.106) 4-(6,7-dichloro-1,2,3,4-tetrahydro-isoquinolin-2-yl-methyl)-aniline

(XV.107) 4-(6,7-dimethyl-1,2,3,4-tetrahydro-isoquinolin-2-yl-methyl)-aniline

(XV.108) 4-(6-chloro-1,2,3,4-tetrahydro-isoquinolin-2-yl-methyl)-aniline

(XV.109) 4-(7-chloro-1,2,3,4-tetrahydro-isoquinolin-2-yl-methyl)-aniline

(XV.110) 4-(6-methoxy-1,2,3,4-tetrahydro-isoquinolin-2-yl-methyl)-aniline

- (XV.111) 4-(7-methoxy-1,2,3,4-tetrahydro-isoquinolin-2-yl-methyl)-aniline
- (XV.112) 4-(2,3,4,5-tetrahydro-azepino(4,5-b)pyrazin-3-yl-methyl)-aniline
- (XV.113) 4-(7-amino-2,3,4,5-tetrahydro-azepino(4,5-b)pyrazin-3-yl-methyl)-aniline
- (XV.114) 4-(2-amino-5,6,7,8-tetrahydro-azepino(4,5-d)thiazol-6-yl-methyl)-aniline
- (XV.115) 4-(5,6,7,8-tetrahydro-azepino(4,5-d)thiazol-6-yl-methyl)-aniline
- (XV.116) 4-(4-methyl-piperazin-1-yl)-aniline
- (XV.117) 4-[N-(2-dimethylamino-ethyl)-N-methyl-amino]-aniline
- (XV.118) 4-[N-(3-dimethylamino-propyl)-N-methyl-amino]-aniline
- (XV.119) N-(3-dimethylamino-propyl)-N-methylsulphonyl-p-phenylenediamine
- (XV.120) 4-[(N-dimethylaminocarbonylmethyl-N-methylsulphonyl)-amino]-aniline
- (XV.121) N-(4-aminophenyl)-N-methyl-methanesulphonamide
- (XV.122) 4-(imidazol-4-yl)-aniline
- (XV.123) 4-(tetrazol-5-yl)-aniline
- (XV.124) 3-[N-(2-dimethylamino-ethyl)-N-propionyl-amino]-aniline
- (XV.125) N-(dimethylamino-methylcarbonyl)-N-methyl-p-phenylenediamine

- (XV.126) N-[(2-dimethylamino-ethyl)-carbonyl]-N-methyl-p-phenylenediamine
- (XV.127) N-dimethylaminocarbonylmethyl-N-acetyl-p-phenylenediamine
- (XV.128) N-methylaminocarbonylmethyl-N-methylsulphonyl-p-phenylenediamine
- (XV.129) N-aminocarbonylmethyl-N-methylsulphonyl-p-phenylenediamine
- (XV.130) 4-(imidazolidin-2,4-dion-5-yliden-methyl)-aniline
- (XV.131) 4-(imidazolidin-2,4-dion-5-yl-methyl)-aniline
- (XV.132) 4-(2-oxo-pyrrolidin-1-yl-methyl)-aniline
- (XV.133) N-cyanomethyl-N-methylsulphonyl-p-phenylenediamine
- (XV.134) 4-[2-(imidazol-4-yl)-ethyl]-aniline
- (XV.135) 4-[(4-methyl-piperazin-1-yl)-methyl]-aniline
- (XV.136) 4-[N-(2-(N-benzyl-N-methyl-amino)-ethyl)-N-methylsulphonyl-amino]-aniline
- (XV.137) 4-[N-(3-(N-benzyl-N-methyl-amino)-propyl)-N-methylsulphonyl-amino]-aniline
- (XV.138) N-cyclohexyl-p-phenylenediamine
- (XV.139) 4-(pyridin-4-yl-methyl)-aniline
- (XV.140) 4-(imidazol-1-yl-methyl)-aniline
- (XV.141) 4-benzyl-aniline

(XV.142) N-(3-trifluoracetyl-amino-propyl)-N-methylsulphonyl-p-phenylenediamine

(XV.143) tert. butyl 4-amino-phenylacetate

(XV.144) 4-(imidazol-2-yl)-aniline

(XV.145) 4-(1-methyl-imidazol-2-yl)-aniline

(XV.146) 4-(1-ethyl-imidazol-2-yl)-aniline

(XV.147) 4-(1-benzyl-imidazol-2-yl)-aniline

(XV.148) 4-[N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino]-3-amino-aniline

(XV.149) 4-[N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino]-3-chloro-aniline

(XV.150) 4-[N-(2-dimethylamino-ethyl)-N-acetyl-amino]-3-amino-aniline

(XV.151) 4-[N-(2-dimethylamino-ethyl)-N-acetyl-amino]-3-bromo-aniline

(XV.152) 4-[2-(4-hydroxy-piperidin-1-yl)-ethyl-amino]-aniline

(XV.153) N-(2-dimethylamino-ethyl)-N-ethylsulphonyl-p-phenylenediamine

(XV.154) N-(2-dimethylamino-ethyl)-N-propylsulphonyl-p-phenylenediamine

(XV.155) N-(2-dimethylamino-ethyl)-N-isopropylsulphonyl-p-phenylenediamine

(XV.156) N-(2-dimethylamino-ethyl)-N-butylsulphonyl-p-phenylenediamine

(XV.157) N-(2-dimethylamino-ethyl)-N-benzylsulphonyl-p-phenylenediamine

(XV.158) N-(2-dimethylamino-ethyl)-N-phenylsulphonyl-p-phenylenediamine

(XV.159) 4-((3-hydroxy-pyrrolidin-1-yl)-methyl)-aniline

(XV.160) 4-[N-(2-dimethylamino-ethyl)-N-(furan-2-carbonyl)-amino]-aniline

(XV.161) 4-[N-(2-dimethylamino-ethyl)-N-(2-methoxy-benzoyl)-amino]-aniline

(XV.162) 4-[N-(2-dimethylamino-ethyl)-N-(pyridine-3-carbonyl)-amino]-aniline

(XV.163) 4-[N-(2-dimethylamino-ethyl)-N-(phenyl-acetyl)-amino]-aniline

(XV.164) N-(piperidin-1-yl-methylcarbonyl)-N-methyl-p-phenylenediamine

(XV.165) N-(morpholin-4-yl-methylcarbonyl)-N-methyl-p-phenylenediamine

(XV.166) N-[(4-benzyl-piperazin-1-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.167) N-(pyrrolidin-1-yl-methylcarbonyl)-N-methyl-p-phenylenediamine

(XV.168) 4-(5-methyl-imidazol-4-yl)-aniline

(XV.169) N-[(2-dimethylamino-ethyl)-carbonyl]-N-isopropyl-p-phenylenediamine

(XV.170) N-[(2-dimethylamino-ethyl)-carbonyl]-N-benzyl-p-phenylenediamine

(XV.171) N-(N-aminocarbonylmethyl-N-methyl-amino)-methylcarbonyl)-N-methyl-p-phenylenediamine

(XV.172) N-[(N-benzyl-N-methyl-amino)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.173) N-[di-(2-methoxyethyl)-amino-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.174) N-[(2-(4-tert.butoxycarbonyl-piperazin-1-yl)-ethyl)-carbonyl]-N-methyl-p-phenylenediamine

(XV.175) N-[(2-(piperidin-1-yl)-ethyl)-carbonyl]-N-methyl-p-phenylenediamine

(XV.176) N-[(2-(N-benzyl-N-methyl-amino)-ethyl)-carbonyl]-N-methyl-p-phenylenediamine

(XV.177) N-(dimethylaminomethylcarbonyl)-N-isopropyl-p-phenylenediamine

(XV.178) N-(piperidin-1-yl-methylcarbonyl)-N-isopropyl-p-phenylenediamine

(XV.179) N-[(4-tert.butoxycarbonyl-piperazin-1-yl)-methylcarbonyl]-N-isopropyl-p-phenylenediamine

(XV.180) N-[(N-benzyl-N-methyl-amino)-methylcarbonyl]-N-benzyl-p-phenylenediamine

(XV.181) N-(dimethylaminomethylcarbonyl)-N-benzyl-p-phenylenediamine

(XV.182) N-(piperidin-1-yl-methylcarbonyl)-N-benzyl-p-phenylenediamine

(XV.183) 4-(1,2,4-triazol-1-yl-methyl)-aniline

(XV.184) 4-(1,2,3-triazol-2-yl-methyl)-aniline

(XV.185) 4-(1,2,3-triazol-1-yl-methyl)-aniline

- (XV.186) 4-[(N-ethoxycarbonylmethyl-N-methyl-amino)-methyl]-aniline
- (XV.187) 4-[(N-aminocarbonylmethyl-N-methyl-amino)-methyl]-aniline
- (XV.188) 4-(azetidin-1-yl-methyl)-aniline
- (XV.189) 4-[(di-(2-methoxy-ethyl)-amino)-methyl]-aniline
- (XV.190) 4-[(N-(2-(2-methoxy-ethoxy)-ethyl)-N-methyl-amino)-methyl]-aniline
- (XV.191) 4-[N-(N-tert.butoxycarbonyl-3-amino-propyl)-N-methyl-aminomethyl]-aniline
- (XV.192) 4-[(N-(methylcarbamoyl-methyl)-N-methyl-amino)-methyl]-aniline
- (XV.193) 4-[(N-(dimethylcarbamoyl-methyl)-N-methyl-amino)-methyl]-aniline
- (XV.194) 4-[(N-propyl-N-methyl-amino)-methyl]-aniline
- (XV.195) 4-[(N-(2-dimethylamino-ethyl)-N-methyl-amino)-methyl]-aniline
- (XV.196) 4-[(N-(3-dimethylamino-propyl)-N-methyl-amino)-methyl]-aniline
- (XV.197) 4-[(N-(2-methoxy-ethyl)-N-methyl-amino)-methyl]-aniline
- (XV.198) 4-[(N-(2-hydroxy-ethyl)-N-methyl-amino)-methyl]-aniline
- (XV.199) 4-[(N-(dioxolan-2-yl-methyl)-N-methyl-amino)-methyl]-aniline
- (XV.200) 4-(3-oxo-piperazin-1-yl-methyl)-aniline
- (XV.201) N-[di-(2-hydroxyethyl)-amino-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.202) N-[(N-(2-methoxyethyl)-N-methyl-amino)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.203) N-[(N-(2-dimethylamino-ethyl)-N-methyl-amino)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.204) N-[(4-methyl-piperazin-1-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.205) N-[(imidazol-1-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.206) N-[(phthalimido-2-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.207) 4-(piperidin-1-yl-methyl)-aniline

The following compounds are prepared analogously to Example XV:

(XV.208) N-(dimethylcarbamoylmethyl)-p-phenylenediamine

(XV.209) di-(2-hydroxyethyl)-aminomethyl-aniline

(XV.210) 4-[N-(2-dimethylamino-ethyl)-N-(methoxyacetyl)-amino]-aniline

(XV.211) 4-[N-(2-dimethylamino-ethyl)-N-(3,4-dimethoxy-benzoyl)-amino]-aniline

(XV.212) 4-[N-(2-(N-benzyl-N-methyl-amino)-ethyl)-N-propionyl-amino]-aniline

(XV.213) 4-[N-(2-dimethylamino-ethyl)-N-(pyridine-4-carbonyl)-amino]-aniline

(XV.214) 4-[N-(2-(N-benzyl-N-methyl-amino)-ethyl)-N-acetyl-amino]-aniline

(XV.215) N-(dimethylaminomethylcarbonyl)-N-methyl-3-methoxy-p-phenylenediamine

(XV.216) N-(dimethylaminomethylcarbonyl)-N-methyl-3-cyano-p-phenylenediamine

(XV.217) 3-(pyridin-4-yl-methyl)-aniline

(XV.218) 4-amino-N-(2-dimethylamino-ethyl)-benzamide

(XV.219) 4-amino-N-(2-dimethylamino-ethyl)-N-methyl-benzamide

(XV.220) 4-amino-N-(3-dimethylamino-propyl)-benzamide

(XV.221) 4-amino-N-(3-dimethylamino-propyl)-N-methyl-benzamide

(XV.222) 4-amino-N-(2-dimethylamino-ethyl)-N-ethyl-benzamide

(XV.223) 4-amino-N-(2-(tert-butyloxycarbonyl-methylamino-ethyl)-N-ethyl-benzamide

(XV.224) 4-amino-N,N-bis-(2-diethylamino-ethyl)-benzamide

(XV.225) 4-amino-N-(2-tert-butyloxycarbonyl-amino-ethyl)-benzamide.

(XV.226) 3-amino-N-(2-dimethylamino-ethyl)-benzamide

(XV.227) 3-amino-N-(2-dimethylamino-ethyl)-N-methyl-benzamide

(XV.228) 3-amino-N-(3-dimethylamino-propyl)-benzamide

(XV.229) 3-amino-N-(3-dimethylamino-propyl)-N-methyl-benzamide

(XV.230) 5-amino-2-N-(dimethylamino-methyl)-carbamoyl-furan

- (XV.231) 4-(4-methyl-piperazin-1-yl-carbonyl)-aniline
- (XV.232) 4-(piperidin-1-yl-carbonyl)-aniline
- (XV.233) 4-amino-N-cyclohexyl-N-methyl-benzamide
- (XV.234) 4-amino-N-isopropyl-benzamide
- (XV.235) 4-(2,3,4,5-tetrahydro-1(H)-benzo[d]azepin-3-yl-carbonyl)-aniline
- (XV.236) 4-(4-hydroxy-piperidin-1-yl-carbonyl)-aniline
- (XV.237) 4-(4-tert-butyloxycarbonyl-piperazin-1-yl-carbonyl)-aniline
- (XV.238) 4-(4-tert-butyloxycarbonyl-[1,4]diazepan-1-yl-carbonyl)-aniline
- (XV.239) 4-amino-1-methyl-2-[N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl]-pyrrole
- (XV.240) 4-amino-1-methyl-2-[(4-methyl-piperazin-1-yl)-carbonyl]-pyrrole
- (XV.241) N-(2-dimethylamino-ethyl)-N-acetyl-2,5-diamino-pyridine
- (XV.242) 4-(1-(2-dimethylamino-ethyl)-imidazol-2-yl)-aniline
- (XV.243) N-[(2-(4-methyl-piperazin-1-yl)-ethyl)-carbonyl]-N-methyl-p-phenylenediamine
- (XV.244) N-[(2-dimethylamino-ethyl)-carbonyl]-N-methyl-p-phenylenediamine
- (XV.245) *trans*-N-dimethylaminomethylcarbonyl-N'-(tert.-butoxycarbonyl)-cyclohexyl-1,4-diamine

(XV.246) 4-(2-dimethylamino-ethoxy)-aniline

(XV.247) 4-[(4-dimethylamino-piperidin-1-yl)-methyl]-aniline

(XV.248) N-[(4-tert.butoxycarbonyl-piperazin-1-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.249) N-hydroxymethylcarbonyl-N-methyl-p-phenylenediamine

(XV.250) 4-[N-(N-tert.butoxycarbonyl-3-aminopropyl)-N-methyl-aminomethyl]-aniline

(XV.251) N-[(4-methyl-homopiperazin-1-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.252) N-[(4-ethyl-piperazin-1-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.253) N-[(4-methyl-piperazin-1-yl)-methylcarbonyl]-p-phenylenediamine

(XV.254) N-[(1-methyl-piperidin-4-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine

(XV.255) 4-(N-(3-dimethylamino-propyl)-N-methyl-carbamoylmethyl)-aniline

(XV.256) 4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoylmethyl)-aniline

(XV.257) 4-[(4-methyl-piperazin-1-yl)-carbonylmethyl]-aniline

(XV.258) N-(4-dimethylaminobutylcarbonyl)-N-methyl-p-phenylenediamine

(XV.259) N-dimethylaminomethylcarbonyl-N'-(tert.-butoxycarbonyl)-2,3-dimethyl-p-phenylenediamine

(XV.260) N-[(4-methyl-piperazin-1-yl)-methylcarbonyl]-N'-(tert.-butoxycarbonyl)-2,3-dimethyl-p-phenylenediamine

(XV.261) N-[(4-methyl-piperazin-1-yl)-aminocarbonyl]-N-methyl-p-phenylenediamine

(XV.262) N-[(4-methyl-piperazin-1-yl)-methylcarbonyl]-N-methyl-m-phenylenediamine

(XV.263) N-[(4-methyl-piperazin-1-yl)-methylcarbonyl]-m-phenylenediamine

(XV.264) N-dimethylaminomethylcarbonyl-N-methyl-m-phenylenediamine

(XV.265) N-[(1-methyl-piperidin-4-yl)-aminocarbonyl]-N-methyl-p-phenylenediamine

(XV.266) N-[(3-dimethylamino-propyl)-carbonyl]-N-methyl-p-phenylenediamine

(XV.267) N-[(4-methyl-piperazin-1-yl)-carbonyl]-N-methyl-p-phenylenediamine

(XV.268) N-[N-(3-dimethylamino-propyl)-aminocarbonyl]-N-methyl-p-phenylenediamine

(XV.269) N-(pyridin-4-yl-methylaminocarbonyl)-N-methyl-p-phenylenediamine

(XV.270) N-(1-methyl-piperidin-4-oxy-carbonyl)-N-methyl-p-phenylenediamine

Example XVI:

trans-N-dimethylaminomethylcarbonyl-cyclohexyl-1,4-diamine-trifluoroacetate
400 mg of *trans*-N-dimethylaminomethylcarbonyl-N'-(tert.-butoxycarbonyl)-cyclohexyl-1,4-diamine (educt XV.245) are dissolved in 12 ml methylene chloride and 5.0 ml of trifluoroacetic acid are added. The mixture is stirred for

0.5 hours at ambient temperature, evaporated down, combined with toluene and concentrated again in the rotary evaporator.

Yield: 420 mg (100% of theory),

$C_{10}H_{21}N_3O$

Mass spectrum: $m/z = 200 [M+H]^+$

The following compounds are prepared analogously to Example XVI:

(XVI.1) N-dimethylaminomethylcarbonyl-2,3-dimethyl-p-phenylenediamine

(XVI.2) N-[(4-methyl-piperazin-1-yl)-methylcarbonyl]-2,3-dimethyl-p-phenylenediamine

Example XVII:

cis-N-dimethylaminomethylcarbonyl-N-methyl-cyclohexyl-1,4-diamin-trifluoroacetate

5.0 g of N-dimethylaminomethylcarbonyl-N-methyl-p-phenylenediamine (educt XV.125) are dissolved in 250 ml glacial acetic acid and 500 mg of Nishimura catalyst (Rh(III)-/Pt(IV)-oxide) are added. The mixture is hydrogenated for 9 hours at ambient temperature under 50 psi, evaporated down, combined with water and then neutralised with sodium hydrogen carbonate solution. After evaporation the residue is taken up in methylene chloride, filtered, dried over sodium sulphate and finally the solvent is eliminated.

Yield: 1.5 g (29% of theory),

$C_{11}H_{23}N_3O$

Mass spectrum: $m/z = 213 [M]^+$

Example XVIII:

N-acetyl-4-(2-diethylamino-ethyl-sulphonyl)-aniline

9.0 g of 4-acetamido-phenylsulphinic acid are dissolved in 10 ml of water and 45 ml of 1N sodium hydroxide solution and 9.47 g of 2-chlorotriethylamine

hydrochloride are added. The mixture is stirred for 5 hours at reflux temperature. After cooling, sodium hydroxide solution is added until an alkaline reaction is obtained, the mixture is extracted with ethyl acetate, dried over magnesium sulphate and finally the solvent is eliminated.

Yield: 9.85 g (73% of theory),

$C_{14}H_{22}N_2O_3S$

Mass spectrum: $m/z = 298 [M]^+$

Example XIX:

4-(2-diethylamino-ethyl-sulphonyl)-aniline

9.85 g of N-acetyl-4-(2-diethylamino-ethyl-sulphonyl)-aniline (educt XVII) are dissolved in 25 ml of ethanol and 100 ml of 3N hydrochloric acid are added.

The mixture is stirred for 1 hour at reflux temperature. After cooling it is neutralised, extracted three times with methylene chloride, dried over magnesium sulphate and finally the solvent is eliminated.

Yield: 5.75 g (68% of theory),

$C_{12}H_{20}N_2O_2S$

Mass spectrum: $m/z = 257 [M+H]^+$

Example XX:

3-Z-[1-(4-(dimethylcarbamoylmethylamino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

6.0 g of 1-acetyl-3-(1-ethoxy-1-phenylmethylen)-6-chloro-2-indolinone (educt IX) and 3.9 g of N-(dimethylcarbamoylmethyl)-p-phenylenediamine (educt XV.208) are dissolved in 50 ml of dimethylformamide and stirred for 4.5 hours at 120°C. After cooling water is added, the precipitate formed is suction filtered and washed with methanol. The product is purified through a silica gel column with methylene chloride/methanol (100:1) as eluant and finally recrystallised from ether.

Yield: 4.4 g (49% of theory),

M.p. 208-211 °C

$C_{27}H_{25}ClN_4O_3$

Mass spectrum: $m/z = 487/489$ $[M-H]^-$

Example XXI:

3-Z-[1-(4-(N-(dimethyl-carbamoyl-methyl)-N-(2-bromo-ethyl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

500 mg of 3-Z-[1-(4-(dimethylcarbamoylmethylamino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone (educt XX) and 430 mg sodium hydrogen carbonate are placed in 10 ml of methylene chloride and 190 mg of 3-bromopropionic acid chloride are slowly added. The mixture is stirred for 1 hour at ambient temperature. Then it is filtered and the filtrate is evaporated down. The precipitate formed is recrystallised from methanol.

Yield: 270 mg (42 % of theory)

R_f value: 0.50 (silica gel, methylene chloride/methanol = 9:1)

$C_{30}H_{28}BrClN_4O_4$

Mass spectrum: $m/z = 621/623/625$ $[M-H]^-$

The following compound is prepared analogously to Example XXI:

(XXI.1) 3-Z-[1-(4-(N-(dimethyl-carbamoyl-methyl)-N-(bromo-acetyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

from 3-Z-[1-(4-(dimethylcarbamoylmethylamino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone (educt XX) and bromoacetylchloride

Preparation of the end compounds:

Example 1.0

3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

0.3 g of 1-acetyl-3-(1-ethoxy-1-phenylmethylene)-6-chloro-2-indolinone (educt IX) and 0.5 g of N-(2-dimethylamino-ethyl)-N-methylsulphonyl-p-phenylenediamine (educt XV.2) are dissolved in 20 ml of dimethylformamide and stirred for 3 hours at 120°C. After cooling 0.8 ml of piperidine are added and the mixture is stirred for a further hour at ambient temperature. The solvent is eliminated and the residue is purified through a silica gel column with methylene chloride/methanol (15:1) as eluant.

Yield: 0.2 g (40% of theory),

R_f value: 0.5 (silica gel, methylene chloride/methanol = 9:1)

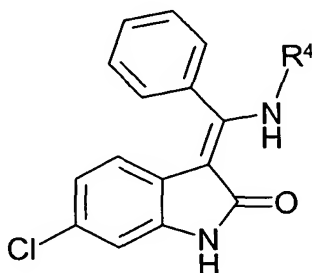
M.p. 237-239 °C

C₂₆H₂₇ClN₄O₃S

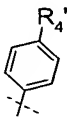
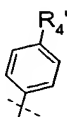
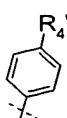
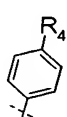
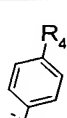
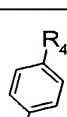
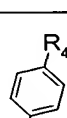
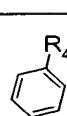
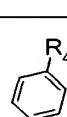
Mass spectrum: m/z = 511/513 [M+H]⁺

The following compounds of general formula I-1 are prepared analogously to

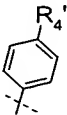
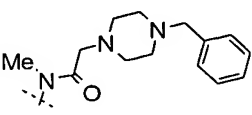
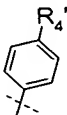
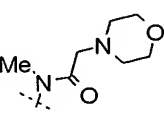
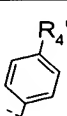
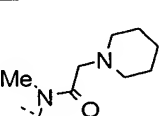
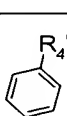
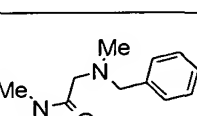
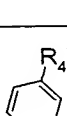
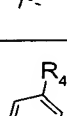
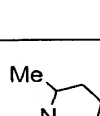
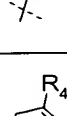
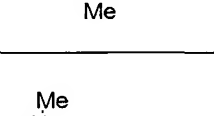
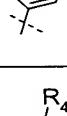
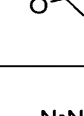
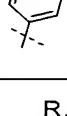
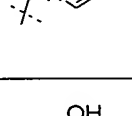
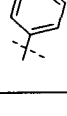
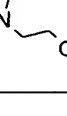
Example 1.0 :



(I-1)

Exam ple	R ⁴	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
1.1		-N(Me)-(CO)-CH ₂ - NMe ₂	XV. 125	C ₂₆ H ₂₅ ClN ₄ O ₂	461/463 [M+H] ⁺	237- 239	0.55 (A)
1.2		-COOEt	-	C ₂₄ H ₁₉ ClN ₂ O ₃	417/419 [M-H] ⁻	266- 268	0.50 (B)
1.3		-N(SO ₂ Me)-(CH ₂)- (CO)-NMe ₂	XV. 120	C ₂₆ H ₂₅ ClN ₄ O ₄ S	523/525 [M-H] ⁻	254	0.50 (C)
1.4		-N(SO ₂ Me)-CH ₃	XV. 121	C ₂₃ H ₂₀ ClN ₃ O ₃ S	452/454 [M-H] ⁻	276- 278	0.50 (C)
1.5		-N(COMe)-CH ₃	-	C ₂₄ H ₂₀ ClN ₃ O ₂	416/418 [M-H] ⁻	308 (decom p.)	0.50 (C)
1.6		-N(SO ₂ Et)-(CH ₂) ₂ - NMe ₂	XV. 153	C ₂₇ H ₂₉ ClN ₄ O ₃ S	523/525 [M-H] ⁻	220	0.50 (C)
1.7		-N(COEt)-(CH ₂) ₂ - NMe ₂	XV.9	C ₂₈ H ₂₉ ClN ₄ O ₂	487/489 [M-H] ⁻	144	0.50 (C)
1.8		-N(SO ₂ Me)-(CH ₂) ₃ - NMe ₂	XV. 119	C ₂₇ H ₂₉ ClN ₄ O ₃ S	524/526 [M] ⁺	214	0.50 (A)
1.9		-N(CO-nPr)-(CH ₂) ₂ - NMe ₂	XV.10	C ₂₉ H ₃₁ ClN ₄ O ₂	501/503 [M-H] ⁻	218	0.50 (A)

1.10		-N(CO-iPr)-(CH ₂) ₂ -NMe ₂	XV.11	C ₂₉ H ₃₁ ClN ₄ O ₂	501/503 [M-H] ⁻	239	0.50 (A)
1.11		-N(COMe)-(CH ₂) ₂ -NMe ₂	XV.6	C ₂₇ H ₂₇ ClN ₄ O ₂	475/477 [M+H] ⁺	170	0.50 (A)
1.12		-N(COPh)-(CH ₂) ₂ -NMe ₂	XV.8	C ₃₂ H ₂₉ ClN ₄ O ₂	537/539 [M+H] ⁺	215	0.50 (A)
1.13		-N(CO-Bn)-(CH ₂) ₂ -NMe ₂	XV. 163	C ₃₃ H ₃₁ ClN ₄ O ₂	551/553 [M+H] ⁺	233	0.50 (A)
1.14		-N[CO-(3-pyridyl)]-(CH ₂) ₂ -NMe ₂	XV. 162	C ₃₁ H ₂₈ ClN ₅ O ₂	538/540 [M+H] ⁺	134	0.50 (A)
1.15		-N[CO-(2-furanyl)]-(CH ₂) ₂ -NMe ₂	XV. 160	C ₃₀ H ₂₇ ClN ₄ O ₃	527/529 [M+H] ⁺	236	0.50 (A)
1.16		-N[CO-(2-MeO-phenyl)]-(CH ₂) ₂ -NMe ₂	XV. 161	C ₃₃ H ₃₁ ClN ₄ O ₃	567/569 [M+H] ⁺	148	0.50 (A)
1.17		-N(SO ₂ -nPr)-(CH ₂) ₂ -NMe ₂	XV. 154	C ₂₈ H ₃₁ ClN ₄ O ₃ S	537/539 [M-H] ⁻	222	0.50 (C)
1.18		-N(SO ₂ -iPr)-(CH ₂) ₂ -NMe ₂	XV. 155	C ₂₈ H ₃₁ ClN ₄ O ₃ S	537/539 [M-H] ⁻	167	0.50 (C)
1.19		-N(SO ₂ Bn)-(CH ₂) ₂ -NMe ₂	XV. 157	C ₃₂ H ₃₁ ClN ₄ O ₃ S	585/587 [M-H] ⁻	132	0.50 (C)

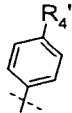
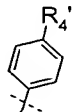
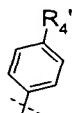
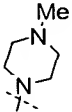
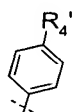
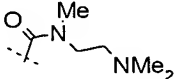
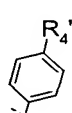
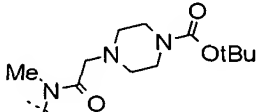
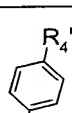
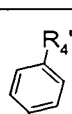
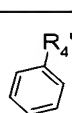
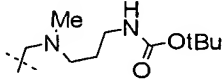
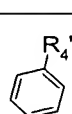
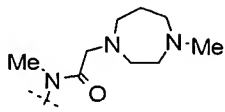
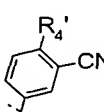
1.20			XV. 166	$C_{35}H_{34}ClN_5O_2$	590/592 [M-H] ⁻	235	0.50 (A)
1.21			XV. 165	$C_{28}H_{27}ClN_4O_3$	501/503 [M-H] ⁻	259	0.50 (A)
1.22			XV. 164	$C_{29}H_{29}ClN_4O_2$	501/503 [M+H] ⁺	235	0.50 (A)
1.23			XV. 172	$C_{26}H_{26}ClN_5O_5S$	535/537 [M-H] ⁻	181	0.50 (A)
1.24		-CH ₂ -NMe ₂	XV.4	$C_{24}H_{22}ClN_3O$	403/405 [M] ⁺	207	0.50 (A)
1.25			XV.1	$C_{29}H_{30}ClN_3O$	470/472 [M-H] ⁻	226	0.50 (A)
1.26			XV. 190	$C_{28}H_{30}ClN_3O_3$	492/494 [M+H] ⁺	140	0.50 (A)
1.27			XV. 185	$C_{24}H_{18}ClN_5O$	450/452 [M+Na] ⁺	230	0.50 (A)
1.28			XV. 209	$C_{26}H_{26}ClN_3O_3$	462/464 [M-H] ⁻	228	0.50 (A)
1.29			XV. 168	$C_{25}H_{19}ClN_4O$	427/429 [M+H] ⁺	290 (decom p.)	0.50 (A)

1.30			XV.13	$C_{29}H_{30}ClN_3O_3$	502/504 [M-H] ⁻	201	0.50 (A)
1.31			XV. 145	$C_{25}H_{19}ClN_4O$	427/429 [M+H] ⁺	279	0.50 (A)
1.32			XV.23	$C_{26}H_{24}ClN_3O_2$	446/448 [M+H] ⁺	245	0.50 (A)
1.33		-CH ₂ -(NBnMe)	XV.78	$C_{30}H_{26}ClN_3O$	502/504 [M+Na] ⁺	168	0.50 (A)
1.34			XV.22	$C_{26}H_{24}ClN_3O$	430/432 [M+H] ⁺	226	0.50 (A)
1.35			XV. 135	$C_{27}H_{27}ClN_4O$	459/461 [M+H] ⁺	228- 230	0.40 (D)
1.36		-N(SO ₂ -nBu)- (CH ₂) ₂ -NMe ₂	XV. 156	$C_{29}H_{33}ClN_4O_3S$	553/555 [M+H] ⁺	185	0.70 (A)
1.37		-N(CO-CH ₂ -OMe)- (CH ₂) ₂ -NMe ₂	XV. 210	$C_{28}H_{29}ClN_4O_3$	505/507 [M+H] ⁺	174	0.40 (A)
1.38		-N[CO-(3,4- dimethoxy-phenyl)]- (CH ₂) ₂ -NMe ₂	XV. 211	$C_{34}H_{33}ClN_4O_4$	597/599 [M+H] ⁺	174	0.50 (A)
1.39			XV. 198	$C_{25}H_{24}ClN_3O_2$	434/436 [M+H] ⁺	208	0.30 (A)

1.50			XV. 174	$C_{34}H_{38}ClN_5O_4$	614/616 [M-H] ⁻	134	0.25 (A)
1.51		-N(COMe)-(CH ₂) ₂ - (NBnMe)	XV. 214	$C_{33}H_{31}ClN_4O_2$	551/553 [M+H] ⁺	195	0.25 (A)
1.52			XV.34	$C_{31}H_{33}ClN_4O_3$	545/547 [M+H] ⁺	225	0.25 (A)
1.53		-N(Me)-(CO)-CH ₂ - NMe ₂	XV. 215	$C_{27}H_{27}ClN_4O_3$	491/493 [M+H] ⁺	238- 241	0.30 (A)
1.54		-CH ₂ -NMe ₂	XV.3	$C_{24}H_{22}ClN_3O$	402/404 [M-H] ⁻	193	0.25 (A)
1.55		-CH ₂ -(4-pyridyl)	XV. 217	$C_{27}H_{20}ClN_3O$	438/440 [M+H] ⁺	243	0.45 (A)
1.56			XII	$C_{26}H_{24}ClN_5O_2$	473/475 [M+H] ⁺	265	0.45 (G)
1.57		-	-	$C_{22}H_{15}ClN_4O$	385/387 [M-H] ⁻	328- 330	0.40 (F)
1.58		-N(COMe)-(CH ₂) ₂ - NMe ₂	XV. 241	$C_{26}H_{26}ClN_5O_2$	476/478 [M+H] ⁺	176- 177	0.60 (H)
1.59		-NH-(CO)-CH ₂ - NMe ₂	XVI	$C_{25}H_{29}ClN_4O_2$	453/455 [M+H] ⁺	n. d.	0.40 (A)

1.60		-N(Me)-(CO)-CH ₂ -NMe ₂	XVII	C ₂₆ H ₃₁ ClN ₄ O ₂	467/469 [M+H] ⁺	257-260	0.20 (F)
1.61			XV. 240	C ₂₆ H ₂₆ ClN ₅ O ₂	476/478 [M+H] ⁺	296-299	0.55 (I)
1.62		-CO-NMe-(CH ₂) ₂ -NMe ₂	XV. 239	C ₂₆ H ₂₈ ClN ₅ O ₂	478/480 [M+H] ⁺	230-232	0.30 (K)
1.63		-N(COMe)-(CH ₂) ₃ -NMe ₂	XV.7	C ₂₈ H ₂₉ ClN ₄ O ₂	489/491 [M+H] ⁺	187	0.20 (A)
1.64		-SO ₂ -(CH ₂) ₂ -NEt ₂	XIX	C ₂₇ H ₂₈ ClN ₃ O ₃ S	510/512 [M+H] ⁺	154-159	0.40 (F)
1.65			XV. 242	C ₂₈ H ₂₆ ClN ₅ O	484/486 [M+H] ⁺	211-216	0.20 (L)
1.66			XV. 232	C ₂₇ H ₂₄ ClN ₃ O ₂	458/460 [M+H] ⁺	270	0.20 (A)
1.67			XV. 237	C ₃₁ H ₃₁ ClN ₄ O ₄	559/561 [M+H] ⁺	255-256	0.20 (A)
1.68			XV. 233	C ₂₉ H ₂₈ ClN ₃ O ₂	486/488 [M+H] ⁺	164	0.20 (A)
1.69		-N(Me)-(CO)-(CH ₂) ₂ -NMe ₂	XV. 126	C ₂₇ H ₂₇ ClN ₄ O ₂	475/477 [M+H] ⁺	219-221	0.20 (A)

1.70		-CH ₂ -NMe-(CH ₂) ₂ -NMe ₂	XV. 195	C ₂₇ H ₂₉ ClN ₄ O	461/463 [M+H] ⁺	151	0.25 (A)
1.71			XV. 243	C ₃₀ H ₃₂ ClN ₅ O ₂	528/520 [M-H] ⁻	204- 208	0.25 (A)
1.72		-O-(CH ₂) ₂ -NMe ₂	XV. 246	C ₂₅ H ₂₄ ClN ₃ O ₂	532/534 [M-H] ⁻	212- 214	0.25 (A)
1.73			XV. 247	C ₂₉ H ₃₁ ClN ₄ O	485/487 [M-H] ⁻	198	0.25 (A)
1.74		-CH ₂ -COOMe	-	C ₂₄ H ₁₉ ClN ₂ O ₃	417/419 [M-H] ⁻	192	0.25 (A)
1.75		-COOMe	-	C ₂₃ H ₁₇ ClN ₂ O ₃	403/405 [M-H] ⁻	209	n. d.
1.76		-(CH ₂) ₂ -NMe ₂	XV.5	C ₂₅ H ₂₄ ClN ₃ O	416/418 [M-H] ⁻	217	0.25 (A)
1.77			XV.12	C ₂₇ H ₂₆ ClN ₃ O ₃	474/476 [M-H] ⁻	203	0.25 (A)
1.78		-N(iPr)-(CO)-CH ₂ -NMe ₂	XV. 177	C ₂₈ H ₂₉ ClN ₄ O ₂	487/489 [M-H] ⁻	216	0.25 (A)
1.79			XV. 179	C ₃₅ H ₄₀ ClN ₅ O ₄	628/630 [M-H] ⁻	164	0.25 (A)

1.80		$-\text{CH}_2-\text{NEt}_2$	XV.66	$\text{C}_{26}\text{H}_{26}\text{ClN}_3\text{O}$	430/432 [M-H] ⁻	244	0.25 (A)
1.81		$-\text{CH}_2-(\text{NMePr})$	XV. 194	$\text{C}_{26}\text{H}_{26}\text{ClN}_3\text{O}$	430/432 [M-H] ⁻	188	0.25 (A)
1.82			XV. 116	$\text{C}_{26}\text{H}_{25}\text{ClN}_4\text{O}$	443/445 [M-H] ⁻	295	0.25 (A)
1.83			XV. 219	$\text{C}_{27}\text{H}_{27}\text{ClN}_4\text{O}_2$	473/475 [M-H] ⁻	148	0.25 (A)
1.84			XV. 248	$\text{C}_{33}\text{H}_{36}\text{ClN}_5\text{O}_4$	602/604 [M+H] ⁺	199	0.25 (A)
1.85		$-\text{N}(\text{Me})-(\text{CO})-\text{CH}_2-\text{OH}$	XV. 249	$\text{C}_{24}\text{H}_{20}\text{ClN}_3\text{O}_3$	432/434 [M-H] ⁻	250	0.25 (A)
1.86		$-\text{N}(\text{Me})-(\text{CH}_2)_2-\text{NMe}_2$	XV. 117	$\text{C}_{26}\text{H}_{27}\text{ClN}_4\text{O}$	445/447 [M-H] ⁻	238	n. d.
1.87			XV. 250	$\text{C}_{31}\text{H}_{35}\text{ClN}_4\text{O}_3$	545/547 [M-H] ⁻	148	0.25 (A)
1.88			XV. 251	$\text{C}_{30}\text{H}_{32}\text{ClN}_5\text{O}_2$	528/530 [M-H] ⁻	223	0.25 (A)
1.89		$-\text{N}(\text{Me})-(\text{CO})-\text{CH}_2-\text{NMe}_2$	XV. 216	$\text{C}_{27}\text{H}_{24}\text{ClN}_5\text{O}_2$	486/488 [M+H] ⁺	226- 228	0.40 (A)

1.90			XV. 252	$C_{30}H_{32}ClN_5O_2$	528/530 [M-H] ⁻	255- 257	0.35 (F)
1.91			XV. 254	$C_{30}H_{31}ClN_4O_2$	515/517 [M+H] ⁺	280- 283	0.30 (M)
1.92			XV. 253	$C_{28}H_{28}ClN_5O_2$	502/504 [M+H] ⁺	251- 255	0.45 (A)
1.93			XV. 255	$C_{29}H_{31}ClN_4O_2$	503/505 [M+H] ⁺	215- 224	0.30 (F)
1.94			XV. 256	$C_{28}H_{29}ClN_4O_2$	489/491 [M+H] ⁺	150- 158	0.40 (A)
1.95			XV. 257	$C_{28}H_{27}ClN_4O_2$	487/489 [M+H] ⁺	244- 248	0.40 (A)
1.96			XV. 258	$C_{29}H_{31}ClN_4O_2$	503/505 [M+H] ⁺	216- 218	0.80 (N)
1.97		-NH-(CO)-CH ₂ - NMe ₂	XVI.1	$C_{27}H_{27}ClN_4O_2$	475/477 [M+H] ⁺	246- 250	0.50 (F)
1.98			XVI.2	$C_{30}H_{32}ClN_5O_2$	530/532 [M+H] ⁺	271- 275	0.50 (F)
1.99			XV. 261	$C_{28}H_{29}ClN_6O_2$	517/519 [M+H] ⁺	250- 253	0.50 (D)

1.100		-CH ₂ -COOEt	-	C ₂₅ H ₂₁ ClN ₂ O ₃	433/435 [M+H] ⁺	166	0.70 (A)
1.101			XV. 265	C ₂₉ H ₃₀ ClN ₅ O ₂	516/518 [M+H] ⁺	265- 270	0.35 (O)
1.102			XV. 266	C ₂₈ H ₂₉ ClN ₄ O ₂	489/491 [M+H] ⁺	238- 242	0.35 (F)
1.103			XV. 267	C ₂₈ H ₂₈ ClN ₅ O ₂	502/504 [M+H] ⁺	290- 293	0.50 (A)
1.104			XV. 268	C ₂₈ H ₃₀ ClN ₅ O ₂	504/506 [M+H] ⁺	192- 195	0.60 (O)
1.105			XV. 269	C ₂₉ H ₂₄ ClN ₅ O ₂	510/512 [M+H] ⁺	222- 223	0.60 (A)
1.106			XV. 270	C ₂₉ H ₂₉ ClN ₄ O ₃	517/519 [M+H] ⁺	237- 240	0.30 (A)

*Eluant mixtures:

(A): silica gel, methylene chloride/methanol 9:1

(B): silica gel, toluene/ethyl acetate 9:1

(C): silica gel, methylene chloride/methanol 10:1

(D): silica gel, methylene chloride/methanol 5:1

(E): silica gel, methylene chloride/methanol/ammonia 5:1:0.01

(F): silica gel, methylene chloride/methanol/ammonia 9:1:0.1

(G): silica gel, methylene chloride/ethanol 5:1

(H): silica gel, methylene chloride/methanol/ammonia 10:1:0.1

(I): silica gel, methylene chloride/ethanol 15:1

(K): silica gel, methylene chloride/ethanol/ammonia 20:1:0.1

(L): silica gel, methylene chloride/ethanol 10:1

(M): silica gel, methylene chloride/methanol/ammonia 9:1:0.01

(N): aluminium oxide, methylene chloride/ethanol 30:1

(O): aluminium oxide, methylene chloride/ethanol 20:1

Example 2.0

3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-5-nitro-2-indolinone

0.4 g of 1-acetyl-3-(1-ethoxy-1-phenylmethylene)-6-chloro-5-nitro-2-indolinone (educt IX.4) and 0.3 g of N-(2-dimethylamino-ethyl)-N-methylsulphonyl-p-phenylenediamine (educt XV.2) are dissolved in 5 ml of dimethylformamide and stirred at 80°C for 4 hours. After cooling, 1.0 ml piperidine is added and the mixture is stirred for a further 3 hours at ambient temperature. The solvent is eliminated and the residue is purified through a silica gel column with methylene chloride/methanol (9:1) as eluant.

Yield: 0.4 g (79% of theory),

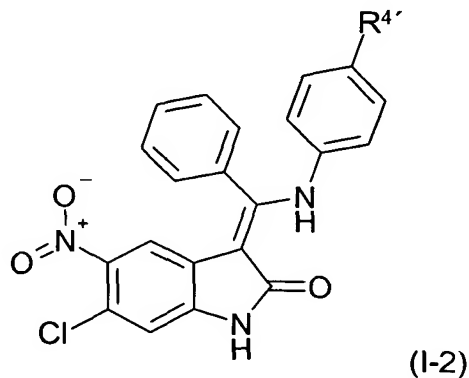
R_f value: 0.5 (silica gel, methylene chloride/methanol = 9:1)

M.p. 224 °C

C₂₆H₂₆ClN₅O₅S

Mass spectrum: m/z = 556/558 [M+H]⁺

The following compounds of general formula I-2 are prepared analogously to Example 2.0:



Exempl e	R ⁴	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
2.1	-N(Me)-(CO)- CH ₂ -NMe ₂	XV.125	C ₂₆ H ₂₄ ClN ₅ O ₄	506/508 [M+H] ⁺	266	0.50 (A)
2.2	-CH ₂ -NMe ₂	XV.4	C ₂₄ H ₂₁ ClN ₄ O ₃	447/449 [M-H] ⁻	260	0.50 (A)
2.3	-N(COMe)- (CH ₂) ₂ -NMe ₂	XV.6	C ₂₇ H ₂₆ ClN ₅ O ₄	520/522 [M+H] ⁺	226	0.50 (A)

*Eluant mixture:

(A): silica gel, methylene chloride/methanol 9:1

Example 3.0

3-Z-[1-(4-(N-(dimethyl-carbamoyl-methyl)-N-(2-pyrrolidin-1-yl-ethyl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

0.3 g of 3-Z-[1-(4-(N-(dimethyl-carbamoyl-methyl)-N-(2-bromo-ethyl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone (educt XXI) and 0.1 ml of pyrrolidine are dissolved in 6 ml of dimethylformamide and stirred for 1.5 hours at ambient temperature. After cooling, 1.1 ml of 1N sodium hydroxide solution are added and the mixture is stirred for another hour at ambient temperature. Water is added, the precipitate formed is suction filtered and purified through a silica gel column with a gradient of methylene chloride and methanol/ammonia as eluant.

Yield: 0.1 g (57% of theory),

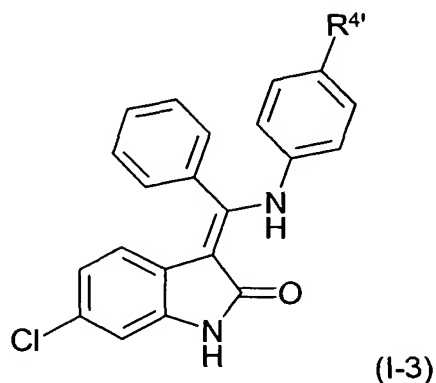
R_f value: 0.20 (silica gel, methylene chloride/methanol/ammonia = 9:1:0.1)

M.p. 224-226 °C

C₃₂H₃₄ClN₅O₃

Mass spectrum: m/z = 570/572 [M-H]⁻

The following compounds of general formula I-3 are prepared analogously to Example 3.0:



Example	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
3.1		XXI.1	C ₃₁ H ₃₂ ClN ₅ O ₃	556/558 [M-H] ⁻	115- 117	0.30 (A)
3.2		XXI	C ₃₀ H ₃₂ ClN ₅ O ₃	546/548 [M+H] ⁺	226	0.25(A)
3.3		XXI.1	C ₂₉ H ₃₀ ClN ₅ O ₃	532/534 [M+H] ⁺	276- 279	0.25 (A)

*Eluant mixture:

(A): silica gel, methylene chloride/methanol/ammonia 9:1:0.1

Example 4.0

3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(3-iodo-phenyl)-methylene]-6-chloro-2-indolinone

0.9 g of 1-acetyl-3-(1-methoxy-1-(3-iodo-phenyl)-methylene)-6-chloro-2-indolinone (educt VIII) and 0.5 g of N-methyl-N-methylsulphonyl-p-phenylenediamine (educt XV.121) are dissolved in 10 ml of dimethylformamide and stirred at 120°C for 3 hours. After cooling, 1.5 ml of

piperidine are added and the mixture is stirred for another hour at ambient temperature. Water is added, the precipitate obtained is suction filtered, washed with a little water, methanol and ether and finally dried in vacuo at 100°C.

Yield: 0.9 g (74% of theory),

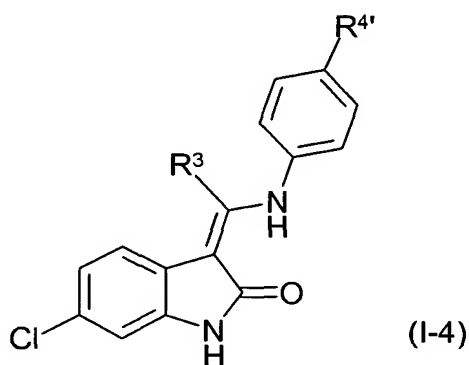
R_f value: 0.6 (silica gel, methylene chloride/methanol = 9:1)

M.p. 292-294 °C

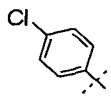
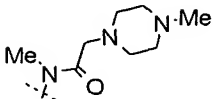
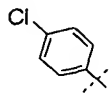
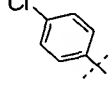
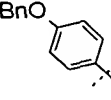
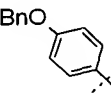
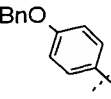
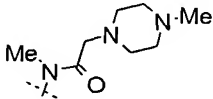
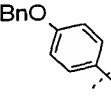
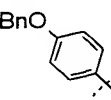
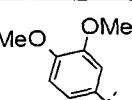
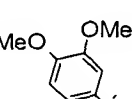
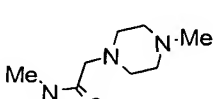
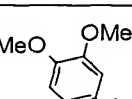
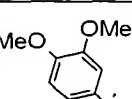
C₂₃H₁₉ClIN₃O₃S

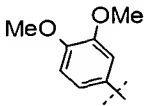
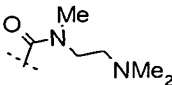
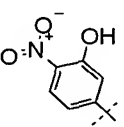
Mass spectrum: m/z = 578/580 [M-H]⁻

The following compounds of general formula I-4 are prepared analogously to Example 4.0 :



Examp le	R ³	R ^{4'}	educts	empirical formula	mass spectrum	m.p. [°C]	R _f value*
4.1		-CH ₂ -NMe ₂	VIII XV.4	C ₂₄ H ₂₁ ClIN ₃ O	529/531 [M+H] ⁺	238- 240	0.30 (A)
4.2		-N(Me)-(CO)- CH ₂ -NMe ₂	VIII.2 XV.125	C ₂₆ H ₂₄ Cl ₂ N ₄ O ₂	495/497 [M+H] ⁺	277- 279	0.20 (B)
4.3		-N(COMe)- (CH ₂) ₂ -NMe ₂	VIII.2 XV.6	C ₂₇ H ₂₆ Cl ₂ N ₄ O ₂	507/509 [M-H] ⁻	241- 243	0.10 (B)

4.4			VIII.2 XV.204	$C_{29}H_{29}Cl_2N_5O_2$	548/550 [M-H] ⁻	266- 268	0.10 (B)
4.5		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.2 XV.7	$C_{28}H_{28}Cl_2N_4O_2$	521/523 [M-H] ⁻	241- 242	0.10 (B)
4.6		-CH ₂ -NMe ₂	VIII.2 XV.4	$C_{24}H_{21}Cl_2N_3O$	438/440 [M+H] ⁺	243- 244	0.10 (B)
4.7		-CH ₂ -NMe ₂	VIII.3 XV.4	$C_{31}H_{28}ClN_3O_2$	510/512 [M+H] ⁺	224- 226	0.30 (B)
4.8		-N(Me)-(CO)- CH ₂ -NMe ₂	VIII.3 XV.125	$C_{33}H_{31}ClN_4O_3$	567/569 [M+H] ⁺	269- 271	0.10 (B)
4.9			VIII.3 XV.204	$C_{36}H_{36}ClN_5O_3$	622/624 [M+H] ⁺	247- 248	0.20 (B)
4.10		-N(COMe)- (CH ₂) ₂ -NMe ₂	VIII.3 XV.6	$C_{34}H_{33}ClN_4O_3$	581/583 [M+H] ⁺	207- 209	0.10 (B)
4.11		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.3 XV.7	$C_{35}H_{35}ClN_4O_3$	595/597 [M+H] ⁺	223- 224	0.10 (B)
4.12		-N(COMe)- (CH ₂) ₂ -NMe ₂	VIII.4 XV.6	$C_{29}H_{31}ClN_4O_4$	533/535 [M-H] ⁻	128- 130	0.75 (C)
4.13			VIII.4 XV.204	$C_{31}H_{34}ClN_5O_4$	574/576 [M-H] ⁻	208- 210	0.65 (C)
4.14		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	VIII.4 XV.2	$C_{28}H_{31}ClN_4O_5S$	569/571 [M-H] ⁻	198- 200	0.75 (C)
4.15		-CH ₂ -NMe ₂	VIII.4 XV.4	$C_{26}H_{26}ClN_3O_3$	462/464 [M-H] ⁻	239- 240	0.70 (C)

4.16			VIII.4 XV.219	$C_{29}H_{31}ClN_4O_4$	533/535 [M-H] ⁻	147- 149	0.70 (C)
4.17		-CH ₂ -NMe ₂	VIII.6 XV.4	$C_{24}H_{21}ClN_4O_4$	465/467 [M+H] ⁺	230 (decom p.)	0.15 (D)

*Eluant mixtures:

(A): silica gel, methylene chloride/methanol 9:1

(B): silica gel, methylene chloride/ethanol 10:1

(C): silica gel, methylene chloride/methanol 4:1

(D): silica gel, methylene chloride/methanol/acetic acid 9:1:0.1

Example 5.0

3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene]-6-chloro-2-indolinone

0.7 g of 1-acetyl-3-(1-chloro-1-(4-(imidazol-1-yl-methyl)-phenyl)-methylene)-6-chloro-2-indolinone (educt X.1), 0.4 g of N-(dimethylamino-methylcarbonyl)-N-methyl-p-phenylenediamine (educt XV.125) and 1.2 ml of triethylamine are dissolved in 10 ml of dimethylformamide and stirred for 15 hours at 60°C. After cooling, 10 ml of methanol and 2 ml concentrated ammonia are added and the mixture is stirred for a further three hours at ambient temperature. Water is added and the mixture is extracted with ethyl acetate. The organic phase is washed three times with water, dried over sodium sulphate and concentrated in the rotary evaporator. The residue is purified through a silica gel column with methylene chloride/methanol/ammonia 10:1:0.1 as eluant. Yield: 0.1 g (5% of theory),

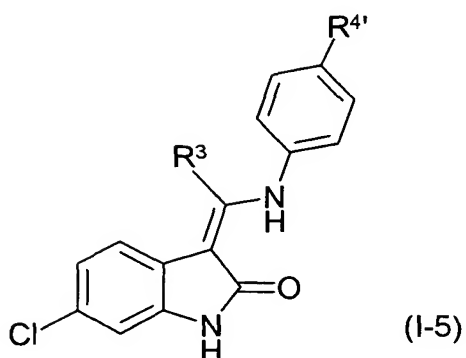
R_f value: not determined

M.p. 268-269 °C

$C_{30}H_{29}ClN_6O_2$

Mass spectrum: m/z = 541/543 [M+H]⁺

The following compounds of general formula I-5 are prepared analogously to Example 5.0:



Example	R ³	R ^{4'}	educts	empirical formula	mass spectrum	m.p. [°C]	R _f value*
5.1		-N(COMe)-(CH ₂) ₂ -NMe ₂	X.1 XV.6	C ₃₁ H ₃₁ ClN ₆ O ₂	555/557 [M+H] ⁺	258- 259	n. d.
5.2			X.1 XV.204	C ₃₃ H ₃₄ ClN ₇ O ₂	594/596 [M-H] ⁻	227	n. d.
5.3		-N(COMe)-(CH ₂) ₃ -NMe ₂	X.1 XV.7	C ₃₂ H ₃₃ ClN ₆ O ₂	567/569 [M-H] ⁻	239- 240	0.20 (A)
5.4			X.2 XV.22	C ₃₀ H ₂₉ ClN ₄ O ₂	511/513 [M-H] ⁻	228- 238	0.30 (B)
5.5		-N(Me)-(CO)-CH ₂ -NMe ₂	X.2 XV.125	C ₃₀ H ₃₀ ClN ₅ O ₃	542/544 [M-H] ⁻	304- 311	0.30 (B)

5.6		-CH ₂ -NMe ₂	X.2 XV.4	C ₂₈ H ₂₇ ClN ₄ O ₂	485/487 [M-H] ⁻	266- 267	0.30 (B)
5.7			X.2 XV.231	C ₃₁ H ₃₀ ClN ₅ O ₃	556/558 [M+H] ⁺	277- 280	0.40 (B)
5.8		-N(COMe)- (CH ₂) ₃ -NMe ₂	X.2 XV.7	C ₃₂ H ₃₄ ClN ₅ O ₃	570/572 [M-H] ⁻	n. d.	0.10 (B)
5.9		-SO ₂ -(CH ₂) ₂ - NEt ₂	X.2 XIX	C ₃₁ H ₃₃ ClN ₄ O ₄ S	591/593 [M-H] ⁻	n. d.	0.40 (B)
5.10		-CH ₂ -NEtMe	X.2 XV.79	C ₂₉ H ₂₉ ClN ₄ O ₂	501/503 [M+H] ⁺	246- 249	0.35 (C)
5.11			X.2 XV.18	C ₃₂ H ₃₃ ClN ₄ O ₄	573/575 [M+H] ⁺	227- 231	0.80 (D)
5.12		-N(Me)-(CO)- CH ₂ -NMe ₂	X XV.125	C ₂₆ H ₂₅ ClN ₆ O ₄	521/523 [M+H] ⁺	254- 256	0.40 (B)
5.13		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	X XV.2	C ₂₆ H ₂₇ ClN ₆ O ₅ S	571/573 [M+H] ⁺	218- 220	0.50 (B)
5.14			X XV.22	C ₂₆ H ₂₄ ClN ₅ O ₃	488/490 [M-H] ⁻	170 (decomp.)	0.30 (B)
5.15			X XV.231	C ₂₇ H ₂₅ ClN ₆ O ₄	531/533 [M-H] ⁻	190- 195	0.30 (E)

5.16		-N(COMe)- (CH ₂) ₂ -NMe ₂	X XV.6	C ₂₇ H ₂₇ ClN ₆ O ₄	533/535 [M-H] ⁻	248- 250	0.30 (F)
5.17		-N(COMe)- (CH ₂) ₃ -NMe ₂	X XV.7	C ₂₈ H ₂₉ ClN ₆ O ₄	547/549 [M-H] ⁻	168- 170	0.30 (G)
5.18		-CH ₂ -NMe ₂	X.3 XV.4	C ₂₄ H ₂₁ ClN ₄ O ₃	447/449 [M-H] ⁻	290- 292	0.30 (H)
5.19		-N(COMe)- (CH ₂) ₂ -NMe ₂	X.3 XV.6	C ₂₇ H ₂₅ ClN ₅ O ₄	518/520 [M-H] ⁻	243- 244	0.35 (I)
5.20			X.3 XV.201	C ₂₉ H ₂₉ ClN ₆ O ₄	559/561 [M-H] ⁻	265- 266	0.25 (I)
5.21		-N(Me)-(CO)- CH ₂ -NMe ₂	X.3 XV.125	C ₂₆ H ₂₄ ClN ₅ O ₄	506/508 [M+H] ⁺	290	0.35 (I)
5.22			X.3 XV.231	C ₂₇ H ₂₄ ClN ₅ O ₄	518/520 [M+H] ⁺	297- 298	0.40 (I)
5.23		-CH ₂ -NMe ₂	X.4 XV.4	C ₂₅ H ₂₁ ClN ₄ O	427/429 [M-H] ⁻	n. d.	0.10 (K)

*Eluant mixtures:

(A): silica gel, methylene chloride/methanol/ammonia 10:1:0.1

(B): silica gel, methylene chloride/methanol/ammonia 9:1:0.1

(C): silica gel, ethyl acetate/methanol/ammonia 7:3:0.1

(D): silica gel, ethyl acetate/ammonia 10:1

(E): silica gel, ethyl acetate/methanol/ammonia 8:2:0.2

(F): silica gel, ethyl acetate/methanol/ammonia 8.5:1.5:0.15

(G): silica gel, methylene chloride/methanol/ammonia 8.5:1.5:0.15

(H): silica gel, methylene chloride/ethanol 5:1

(I): silica gel, methylene chloride/ethanol/ammonia 20:1:0.1

(K): silica gel, methylene chloride/methanol 9:1

Example 6.0

3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-bromo-2-indolinone

1.0 g of 1-acetyl-3-(1-ethoxy-1-phenylmethylene)-6-bromo-2-indolinone (educt IX.1) and 0.7 g of N-(2-dimethylamino-ethyl)-N-methylsulphonyl-p-phenylenediamine (educt XV.2) are dissolved in 4 ml of dimethylformamide and stirred for 2 hours at 120°C. After cooling a little methanol is added and the precipitate formed is suction filtered. Then the residue is suspended in a little ethanol, 3.3 ml of 1 N sodium hydroxide solution are added and the mixture is stirred for another hour at ambient temperature. After this time water is added, the precipitate is suction filtered and washed with water, methanol and ether.

Yield: 0.7 g (50% of theory),

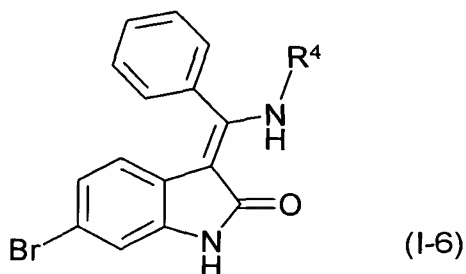
R_f value: 0.35 (silica gel, methylene chloride/methanol = 9:1)

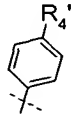
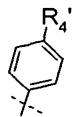
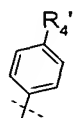
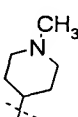
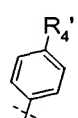
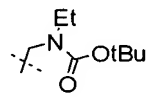
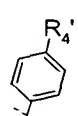

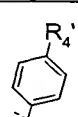
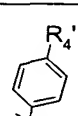
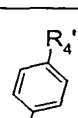
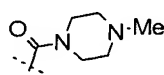
M.p. 204-205 °C

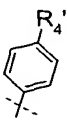
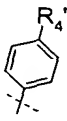
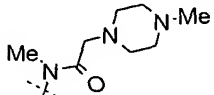
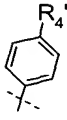
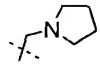
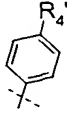
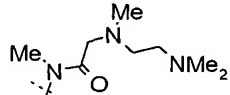
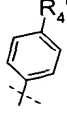
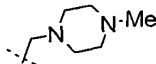
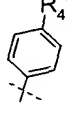
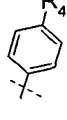
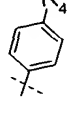
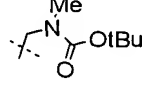
C₂₆H₂₇BrN₄O₃S

Mass spectrum: m/z = 555/557 [M]⁺

The following compounds of general formula I-6 are prepared analogously to Example 6.0:



Example	R ⁴	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
6.1		-N(Me)-(CO)-CH ₂ -NMe ₂	XV.125	C ₂₆ H ₂₅ BrN ₄ O ₂	505/507 [M+H] ⁺	253-256	0.35 (A)
6.2		-CH ₂ -NMe ₂	XV.4	C ₂₄ H ₂₂ BrN ₃ O	448/450 [M+H] ⁺	236-238	0.20 (A)
6.3		-N(COMe)-(CH ₂) ₂ -NMe ₂	XV.6	C ₂₇ H ₂₇ BrN ₄ O ₂	517/519 [M-H] ⁻	147	0.25 (A)
6.4		-	-	C ₂₁ H ₂₂ BrN ₃ O	411/413 [M] ⁺	358	0.20 (A)
6.5			XV.13	C ₂₉ H ₃₀ BrN ₃ O ₃	546/548 [M-H] ⁻	186-188	0.60 (A)
6.6			XV.145	C ₂₅ H ₁₉ BrN ₄ O	469/471 [M-H] ⁻	302-304	0.50 (A)
6.7		-N(SO ₂ Bn)-(CH ₂) ₂ -NMe ₂	XV.157	C ₃₂ H ₃₁ BrN ₄ O ₃ S	629/631 [M-H] ⁻	131-134	0.25 (A)
6.8		-N(SO ₂ nPr)-(CH ₂) ₂ -NMe ₂	XV.153	C ₂₈ H ₃₁ BrN ₄ O ₃ S	581/583 [M-H] ⁻	228-230	0.25 (A)
6.9			XV.231	C ₂₇ H ₂₅ BrN ₄ O ₂	515/517 [M-H] ⁻	268-270	0.25 (A)

6.10		$\text{-N(COMe)-(CH}_2\text{)}_3\text{-NMe}_2$	XV.7	$\text{C}_{28}\text{H}_{29}\text{BrN}_4\text{O}_2$	531/533 [M-H] ⁻	138	0.25 (A)
6.11			XV.204	$\text{C}_{29}\text{H}_{30}\text{BrN}_5\text{O}_2$	560/562 [M+H] ⁺	276- 278	0.40 (A)
6.12			XV.22	$\text{C}_{26}\text{H}_{24}\text{BrN}_3\text{O}$	474/476 [M+H] ⁺	243- 247	0.50 (B)
6.13			XV.203	$\text{C}_{29}\text{H}_{32}\text{BrN}_5\text{O}_2$	562/564 [M+H] ⁺	178	0.60 (C)
6.14			XV.135	$\text{C}_{27}\text{H}_{27}\text{BrN}_4\text{O}$	503/505 [M+H] ⁺	247	0.70 (D)
6.15		$\text{-N(Me)-(CO)-(CH}_2\text{)}_2\text{-NMe}_2$	XV.126	$\text{C}_{27}\text{H}_{27}\text{BrN}_4\text{O}_2$	519/521 [M+H] ⁺	229	0.30 (D)
6.16		$\text{-CH}_2\text{-NMe-(CH}_2\text{)}_2\text{-NMe}_2$	XV.195	$\text{C}_{27}\text{H}_{29}\text{BrN}_4\text{O}$	505/507 [M+H] ⁺	160	0.20 (D)
6.17			XV.18	$\text{C}_{28}\text{H}_{28}\text{BrN}_3\text{O}_3$	532/534 [M-H] ⁻	212- 215	0.55 (E)

(A): silica gel, methylene chloride/methanol 9:1

(B): silica gel, methylene chloride/methanol/ammonia 9:1:0.01

(C): aluminium oxide, methylene chloride/methanol 9:1

(D): silica gel, methylene chloride/methanol/ammonia 5:1:0.1

(E): silica gel, methylene chloride/methanol/ammonia 9:1:0.1

Example 7.03-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone

67 mg of 1-acetyl-3-(1-ethoxy-1-phenylmethylene)-6-cyano-2-indolinone (educt IX.2) and 60 mg of N-(2-dimethylamino-ethyl)-N-methylsulphonyl-p-phenylenediamine (educt XV.2) are dissolved in 5 ml of dimethylformamide and stirred for 1.5 hours at 80°C. After cooling, 2 ml concentrated ammonia solution are added and the mixture is stirred for a further 10 minutes at ambient temperature. After this time water is added, the precipitate is suction filtered, dissolved again in methylene chloride/methanol and dried over sodium sulphate. After removal of the solvent the residue is washed with ether and dried at 80°C.

Yield: 28 mg (26% of theory),

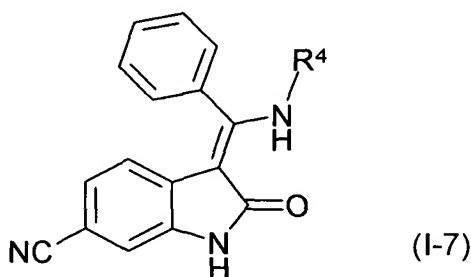
R_f value: 0.15 (silica gel, methylene chloride/methanol = 9:1)

M.p. 270 °C

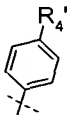
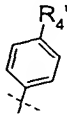
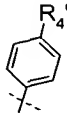
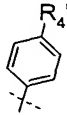
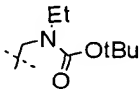
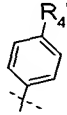
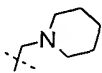
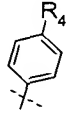
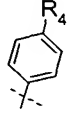
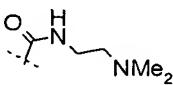
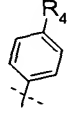
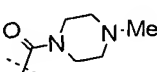
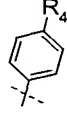
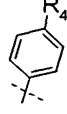
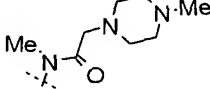
C₂₇H₂₇N₅O₃S

Mass spectrum: m/z = 501 [M]⁺

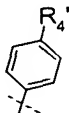
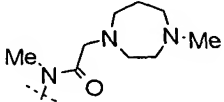
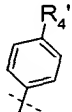
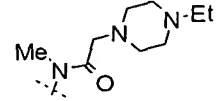
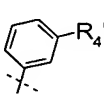
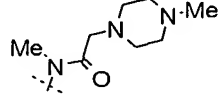
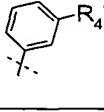
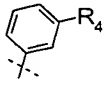
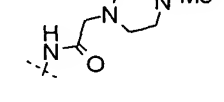
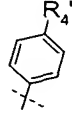
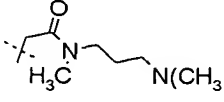
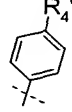
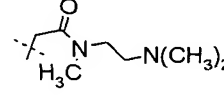
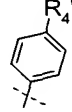
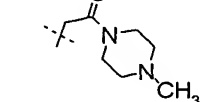
The following compounds of general formula I-7 are prepared analogously to Example 7.0:



Example	R ⁴	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
7.1		-N(Me)-(CO)-CH ₂ -NMe ₂	XV.125	C ₂₇ H ₂₅ N ₅ O ₂	452 [M+H] ⁺	263-266	0.10 (A)

7.2		-CH ₂ -NMe ₂	XV.4	C ₂₅ H ₂₂ N ₄ O	393 [M-H] ⁻	267- 269	0.60 (B)
7.3		-N(COMe)-(CH ₂) ₂ - NMe ₂	XV.6	C ₂₈ H ₂₇ N ₅ O ₂	464 [M-H] ⁻	277- 280	0.40 (A)
7.4		-Br	-	C ₂₂ H ₁₄ BrN ₃ O	414/416 [M-H] ⁻	338- 340	0.30 (A)
7.5			XV.13	C ₃₀ H ₃₀ N ₄ O ₃	493 [M-H] ⁻	201- 204	0.55 (A)
7.6			XV.207	C ₂₈ H ₂₆ N ₄ O	433 [M-H] ⁻	259	0.25 (A)
7.7		-O-(CH ₂) ₂ -NMe ₂	XV.246	C ₂₆ H ₂₄ N ₄ O ₂	423 [M-H] ⁻	256- 258	0.50 (B)
7.8			XV.218	C ₂₇ H ₂₅ N ₄ O ₂	450 [M-H] ⁻	258- 260	0.20 (B)
7.9			XV.231	C ₂₈ H ₂₅ N ₅ O ₂	462 [M-H] ⁻	328- 329	0.75 (B)
7.10		-N(COMe)-(CH ₂) ₃ - NMe ₂	XV.7	C ₂₉ H ₂₉ N ₅ O ₂	478 [M-H] ⁻	262	n. d.
7.11			XV.204	C ₃₀ H ₃₀ N ₆ O ₂	507 [M+H] ⁺	305- 307	0.15 (A)

7.12			XV.22	$C_{27}H_{24}N_4O$	421 [M+H] ⁺	248	0.10 (A)
7.13			XV.203	$C_{30}H_{32}N_6O_2$	509 [M+H] ⁺	218- 220	0.40 (C)
7.14			XV.220	$C_{28}H_{27}N_5O_2$	466 [M+H] ⁺	247- 249	0.10 (B)
7.15			XV.219	$C_{28}H_{27}N_5O_2$	466 [M+H] ⁺	208- 210	0.45 (B)
7.16			XV.221	$C_{29}H_{29}N_5O_2$	480 [M+H] ⁺	264- 267	0.10 (B)
7.17		-N(Me)-(CO)- (CH ₂) ₂ -NMe ₂	XV.126	$C_{28}H_{27}N_5O_2$	466 [M+H] ⁺	274	0.15 (A)
7.18			XV.248	$C_{34}H_{36}N_6O_4$	593 [M+H] ⁺	251- 254	0.30 (A)
7.19		-CH ₂ -NMe ₂	XV.3	$C_{25}H_{22}N_4O$	393 [M-H] ⁻	232	0.35 (A)
7.20			XV.227	$C_{28}H_{27}N_5O_2$	466 [M+H] ⁺	188- 191	0.40 (C)
7.21			XV.165	$C_{29}H_{27}N_5O_3$	494 [M+H] ⁺	301	0.30 (A)

7.22			XV.251	$C_{31}H_{32}N_6O_2$	519 [M-H] ⁻	250	n. d.
7.23			XV.252	$C_{31}H_{32}N_6O_2$	519 [M-H] ⁻	276	0.45 (B)
7.24			XV.262	$C_{30}H_{30}N_6O_2$	507 [M+H] ⁺	199	0.50 (B)
7.25		-N(Me)-(CO)-CH ₂ - NMe ₂	XV.264	$C_{27}H_{25}N_5O_2$	452 [M+H] ⁺	199	0.50 (B)
7.26			XV.263	$C_{29}H_{28}N_6O_2$	493 [M+H] ⁺	196	0.30 (A)
7.27			XV.255	$C_{30}H_{31}N_5O_2$	494 [M+H] ⁺	201	0.45 (C)
7.28			XV.256	$C_{29}H_{29}N_5O_2$	480 [M+H] ⁺	206	0.25 (C)
7.29			XV.257	$C_{29}H_{27}N_5O_2$	478 [M+H] ⁺	256	0.30 (C)

(A): silica gel, methylene chloride/methanol 9:1

(B): silica gel, methylene chloride/methanol 4:1

(C): silica gel, methylene chloride/methanol/ammonia 9:1:0.1

Example 8.03-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone

325 mg of 1-acetyl-3-(1-ethoxy-1-phenylmethylene)-6-fluoro-2-indolinone (educt IX.3) and 310 mg N-(2-dimethylamino-ethyl)-N-methylsulphonyl-p-phenylenediamine (educt XV.2) are dissolved in 2 ml of dimethylformamide and stirred for 4 hours at 120°C. After cooling methanol and water are added, the mixture is extracted with ethyl acetate and the organic phase is concentrated in the rotary evaporator. The residue obtained is purified through a silica gel column with methylene chloride/methanol 9:1 as eluant. The product is suspended in a little ethanol to eliminate the acetyl group, 1.3 ml of 1 N sodium hydroxide solution are added and the mixture is stirred for another hour at ambient temperature. After this time water is added, the precipitate is suction filtered and washed with water, methanol and ether. The residue is dried in vacuo at 100°C.

Yield: 0.3 g (61% of theory),

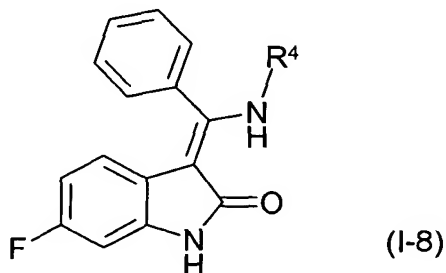
R_f value: 0.25 (silica gel, methylene chloride/methanol/ammonia = 9:1:0.1)

M.p. 259-261 °C

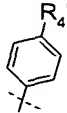

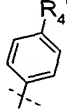
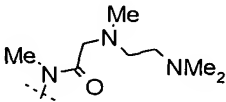
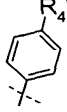
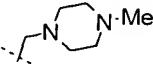
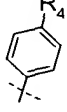
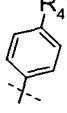
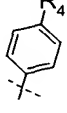

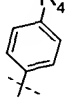
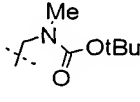
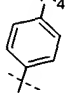

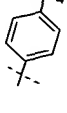
C₂₆H₂₇FN₄O₃S


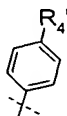
Mass spectrum: m/z = 493 [M-H]⁺

The following compounds of general formula I-8 are prepared analogously to Example 8.0:



Examp le	R ⁴	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
8.1		-N(Me)-(CO)-CH ₂ - NMe ₂	XV.125	C ₂₆ H ₂₅ FN ₄ O ₂	445 [M+H] ⁺	226	0.25 (A)
8.2		-CH ₂ -NMe ₂	XV.4	C ₂₄ H ₂₂ FN ₃ O	386 [M-H] ⁻	229- 232	0.35 (A)
8.3		-N(COMe)-(CH ₂) ₂ - NMe ₂	XV.6	C ₂₇ H ₂₇ FN ₄ O ₂	459 [M+H] ⁺	225- 227	0.25 (A)
8.4			XV.13	C ₂₉ H ₃₀ FN ₃ O ₃	486 [M-H] ⁻	182	0.50 (B)
8.5			XV.145	C ₂₅ H ₁₉ FN ₄ O	411 [M+H] ⁺	290	0.50 (C)
8.6		-N(SO ₂ nPr)- (CH ₂) ₂ -NMe ₂	XV.154	C ₂₈ H ₃₁ FN ₄ O ₃ S	521 [M-H] ⁻	227	0.35 (A)
8.7			XV.231	C ₂₇ H ₂₅ FN ₄ O ₂	457 [M+H] ⁺	118	0.35 (A)
8.8		-N(COMe)-(CH ₂) ₃ - NMe ₂	XV.7	C ₂₈ H ₂₉ FN ₄ O ₂	473 [M+H] ⁺	214	0.25 (A)
8.9			XV.204	C ₂₉ H ₃₀ FN ₅ O ₂	500 [M+H] ⁺	230	0.30 (A)

8.10			XV.22	$C_{26}H_{24}FN_3O$	414 [M+H] ⁺	220	0.25 (A)
8.11			XV.203	$C_{29}H_{32}FN_5O_2$	500 [M-H] ⁻	150	0.25 (D)
8.12			XV.135	$C_{27}H_{27}FN_4O$	443 [M+H] ⁺	198	0.15 (A)
8.13		-N(Me)-(CO)- (CH ₂) ₂ -NMe ₂	XV.126	$C_{27}H_{27}FN_4O_2$	459 [M+H] ⁺	201	0.40 (A)
8.14		-CH ₂ -NMe-(CH ₂) ₂ - NMe ₂	XV.195	$C_{27}H_{29}FN_4O$	445 [M+H] ⁺	141	0.30 (A)
8.15			XV.207	$C_{27}H_{26}FN_3O$	426 [M-H] ⁻	223	0.60 (A)
8.16			XV.18	$C_{28}H_{28}FN_3O_3$	474 [M+H] ⁺	168	0.55 (B)
8.17		-NMe-SO ₂ Me	XV.121	$C_{23}H_{20}FN_3O_3S$	438 [M+H] ⁺	295- 300	0.60 (A)
8.18		-SO ₂ Me	-	$C_{22}H_{17}FN_2O_3S$	409 [M+H] ⁺	255- 260	0.50 (A)
8.19		-N(COMe)-CH ₃	-	$C_{24}H_{20}FN_3O_2$	402 [M+H] ⁺	310- 315	0.45 (A)

8.20		-	-	$C_{21}H_{15}FN_2O$	331 [M+H] ⁺	299- 305	0.60 (A)
8.21		-N(SO ₂ Me)-(CH ₂)- (CO)-NMe ₂	XV.120	$C_{26}H_{25}FN_4O_4S$	509 [M+H] ⁺	270- 274	0.50 (E)

(A): silica gel, methylene chloride/methanol/ammonia 9:1:0.1

(B): silica gel, methylene chloride/methanol/ammonia 10:1:0.1

(C): silica gel, methylene chloride/methanol 9:1

(D): silica gel, methylene chloride/methanol/ammonia 7:1:0.1

(E): aluminium oxide, methylene chloride/methanol 19:1

Example 9.0

3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(4-iodo-phenyl)-methylene]-6-fluoro-2-indolinone

3.5 g of 1-acetyl-3-(1-methoxy-1-(4-iodo-phenyl)-methylene)-6-fluoro-2-indolinone (educt VIII.14) and 1.6 g of 4-(dimethylaminomethyl)-aniline (educt XV.4) are dissolved in 30 ml of dimethylformamide and stirred for 2 hours at 120°C. After cooling the solvent is eliminated, the residue is taken up in 30 ml of methanol and 2 spatula tips of sodium methoxide are added. After a yellow precipitate has formed the solvent is removed by suction filtering, the residue is washed with a little methanol and ether and finally dried in vacuo at 100°C. Yield: 1.9 g (46% of theory),

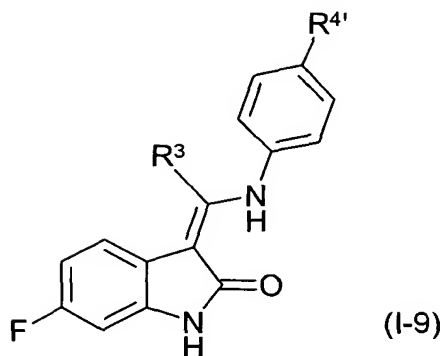
R_f value: 0.3 (silica gel, methylene chloride/methanol = 9:1)

M.p. 243-246 °C

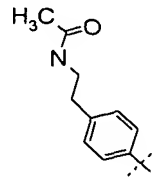
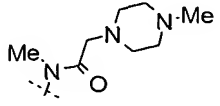
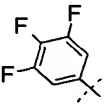
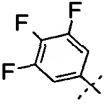
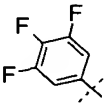
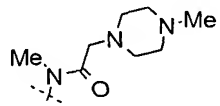
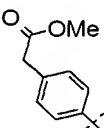
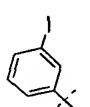
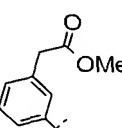
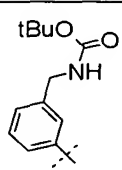
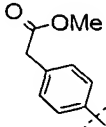
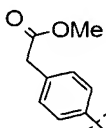
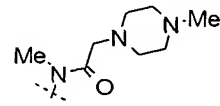
$C_{24}H_{21}FIN_3O$

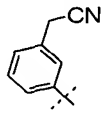
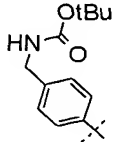
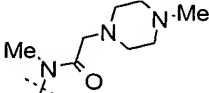
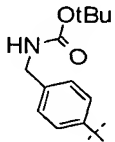
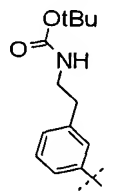
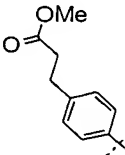
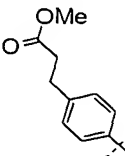
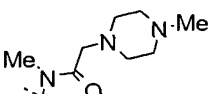
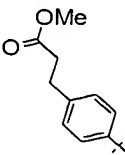
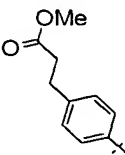
Mass spectrum: m/z = 514 [M+H]⁺

The following compounds of general formula I-9 are prepared analogously to Example 9.0:

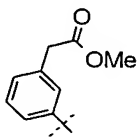
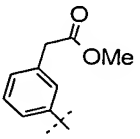
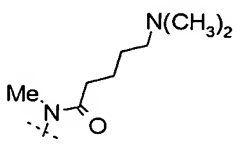
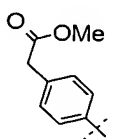
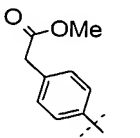

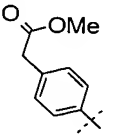
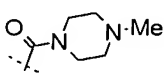
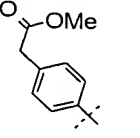
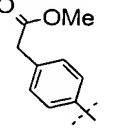
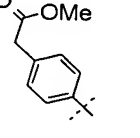
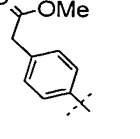


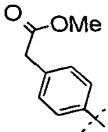
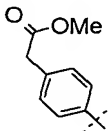
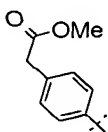
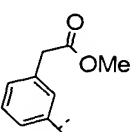
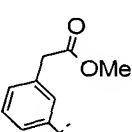
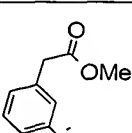

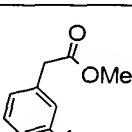
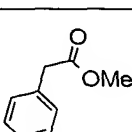
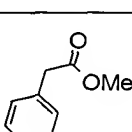
Example	R ³	R ^{4'}	educts	empirical formula	mass spectrum	m.p. [°C]	R _f value*
9.1		-CH ₂ -NMe ₂	VIII.7 XV.4	C ₂₄ H ₂₁ F ₂ N ₃ O	404 [M-H] ⁻	225- 227	0.20 (A)
9.2		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.7 XV.7	C ₂₈ H ₂₈ F ₂ N ₄ O ₂	491 [M+H] ⁺	160- 163	0.20 (A)
9.3			VIII.7 XV.204	C ₂₉ H ₂₉ F ₂ N ₅ O ₂	518 [M+H] ⁺	218- 220	0.40 (A)
9.4		-CH ₂ -NMe ₂	VIII.8 XV.4	C ₂₈ H ₂₉ FN ₄ O ₂	471 [M-H] ⁻	106- 110	0.25 (A)
9.5		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.8 XV.7	C ₃₂ H ₃₆ FN ₅ O ₃	558 [M+H] ⁺	194- 196	0.25 (A)

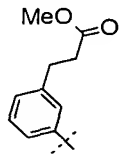
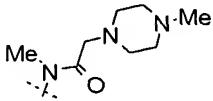
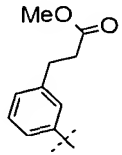
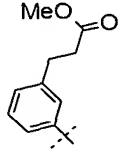
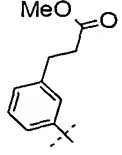
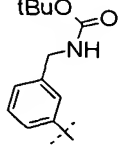
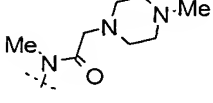
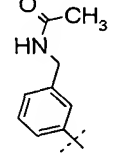
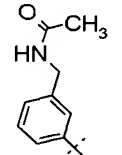
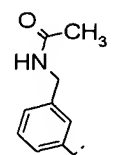
9.6			VIII.8 XV.204	$C_{33}H_{37}FN_6O_3$	583 [M-H] ⁻	238- 240	0.25 (A)
9.7		-CH ₂ -NMe ₂	VIII.9 XV.4	$C_{24}H_{19}F_4N_3O$	440 [M-H] ⁻	267- 269	0.35 (A)
9.8		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.9 XV.7	$C_{28}H_{26}F_4N_4O_2$	527 [M+H] ⁺	210- 212	0.15 (A)
9.9			VIII.9 XV.204	$C_{29}H_{27}F_4N_5O_2$	554 [M+H] ⁺	216- 218	0.20 (A)
9.10		-CH ₂ -NMe ₂	VIII.1 XV.4	$C_{27}H_{26}FN_3O_3$	460 [M+H] ⁺	173- 176	0.30 (A)
9.11		-CH ₂ -NMe ₂	VIII.16 XV.4	$C_{24}H_{21}FIN_3O$	514 [M+H] ⁺	198- 200	0.30 (B)
9.12		-CH ₂ -NMe ₂	VIII.10 XV.4	$C_{27}H_{26}FN_3O_3$	458 [M-H] ⁻	195- 198	0.25 (A)
9.13		-CH ₂ -NMe ₂	VIII.11 XV.4	$C_{30}H_{33}FN_4O_3$	517 [M+H] ⁺	230- 240	0.30 (A)
9.14		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	VIII.1 XV.2	$C_{29}H_{31}FN_4O_5S$	567 [M+H] ⁺	188- 189	0.40 (A)
9.15			VIII.1 XV.204	$C_{32}H_{34}FN_5O_4$	572 [M+H] ⁺	200- 203	0.35 (C)

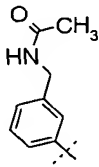
9.16		-CH ₂ -NMe ₂	VIII.12 XV.4	C ₂₆ H ₂₃ FN ₄ O	427 [M+H] ⁺	130- 135	0.25 (A)
9.17			VIII.13 XV.204	C ₃₅ H ₄₁ FN ₆ O ₄	629 [M+H] ⁺	215- 220	0.35 (A)
9.18		-CH ₂ -NMe ₂	VIII.13 XV.4	C ₃₀ H ₃₃ FN ₄ O ₃	517 [M+H] ⁺	186- 190	0.35 (A)
9.19		-CH ₂ -NMe ₂	VIII.20 XV.4	C ₃₁ H ₃₅ FN ₄ O ₃	531 [M+H] ⁺	n. d.	0.40 (A)
9.20		-NMe-(COMe)	VIII.18 -	C ₂₈ H ₂₆ FN ₃ O ₄	488 [M+H] ⁺	166- 170	0.40 (A)
9.21			VIII.18 XV.204	C ₃₃ H ₃₆ FN ₅ O ₄	586 [M+H] ⁺	176- 180	0.30 (A)
9.22		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	VIII.18 XV.2	C ₃₀ H ₃₃ FN ₄ O ₅ S	581 [M+H] ⁺	195- 198	0.45 (A)
9.23		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.18 XV.7	C ₃₂ H ₃₅ FN ₄ O ₄	559 [M+H] ⁺	100- 104	0.50 (A)

9.24			VIII.18 XV.18	$C_{32}H_{34}FN_3O_5$	558 [M-H] ⁻	132- 137	0.80 (D)
9.25			VIII.18 XV.231	$C_{31}H_{31}FN_4O_4$	543 [M+H] ⁺	234- 236	0.60 (A)
9.26			VIII.18 XV.145	$C_{29}H_{25}FN_4O_3$	497 [M+H] ⁺	110- 115	0.40 (A)
9.27		-SO ₂ Me	VIII.18 -	$C_{26}H_{23}FN_2O_5S$	495 [M+H] ⁺	130- 137	0.60 (A)
9.28			VIII.10 XV.204	$C_{32}H_{34}FN_5O_4$	572 [M+H] ⁺	189	0.60 (B)
9.29		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	VIII.10 XV.2	$C_{29}H_{31}FN_4O_5S$	567 [M+H] ⁺	n. d.	0.60 (B)
9.30			VIII.10 XV.231	$C_{30}H_{29}FN_4O_4$	529 [M+H] ⁺	201- 203	0.60 (B)
9.31		-N(Me)-(CO)- CH ₂ -NMe ₂	VIII.10 XV.125	$C_{29}H_{29}FN_4O_4$	517 [M+H] ⁺	126	0.60 (B)
9.32		-N(COMe)- (CH ₂) ₂ -NMe ₂	VIII.10 XV.6	$C_{30}H_{31}FN_4O_4$	531 [M+H] ⁺	179	0.50 (B)

9.33		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.10 XV.7	C ₃₁ H ₃₃ FN ₄ O ₄	545 [M+H] ⁺	123	0.20 (B)
9.34			VIII.10 XV.258	C ₃₂ H ₃₅ FN ₄ O ₄	559 [M+H] ⁺	201	0.20 (B)
9.35		-H	VIII.1 -	C ₂₄ H ₁₉ FN ₂ O ₃	403 [M+H] ⁺	198- 206	0.80 (A)
9.36			VIII.1 XV.145	C ₂₈ H ₂₃ FN ₄ O ₃	483 [M+H] ⁺	223- 226	0.75 (A)
9.37			VIII.1 XV.231	C ₃₀ H ₂₉ FN ₄ O ₄	529 [M+H] ⁺	215- 220	0.30 (A)
9.38		-N(SO ₂ Me)- (CH ₂)-(CO)- NMe ₂	VIII.1 XV.120	C ₂₉ H ₂₉ FN ₄ O ₆ S	581 [M+H] ⁺	227- 230	0.65 (A)
9.39		-N(Me)-(CO)- CH ₂ -NMe ₂	VIII.1 XV.125	C ₂₉ H ₂₉ FN ₄ O ₄	517 [M+H] ⁺	128- 130	0.45 (A)
9.40		-N(COMe)-CH ₃	VIII.1 -	C ₂₇ H ₂₄ FN ₃ O ₄	474 [M+H] ⁺	218- 223	0.40 (A)
9.41		-N(Me)-(CO)- (CH ₂) ₂ -NMe ₂	VIII.1 XV.126	C ₃₀ H ₃₁ FN ₄ O ₄	531 [M+H] ⁺	192- 194	0.40 (A)

9.42		-SO ₂ Me	VIII.1 -	C ₂₅ H ₂₁ FN ₂ O ₅ S	481 [M+H] ⁺	205- 214	0.65 (A)
9.43		-N(Me)-(CO)- (CH ₂) ₃ -NMe ₂	VIII.1 XV.268	C ₃₁ H ₃₃ FN ₄ O ₄	545 [M+H] ⁺	190- 193	0.15 (A)
9.44		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.1 XV.7	C ₃₁ H ₃₃ FN ₄ O ₄	545 [M+H] ⁺	184- 188	0.50 (A)
9.45		-H	VIII.10 -	C ₂₄ H ₁₉ FN ₂ O ₃	403 [M+H] ⁺	114	0.70 (B)
9.46		-SO ₂ Me	VIII.10 -	C ₂₅ H ₂₁ FN ₂ O ₅ S	481 [M+H] ⁺	129	0.60 (B)
9.47			VIII.10 XV.145	C ₂₈ H ₂₃ FN ₄ O ₃	483 [M+H] ⁺	125	0.60 (B)
9.48		-N(SO ₂ Me)- (CH ₂)-(CO)- NMe ₂	VIII.10 XV.120	C ₂₉ H ₂₉ FN ₄ O ₆ S	581 [M+H] ⁺	163	0.60 (B)
9.49		-N(Me)-(CO)- (CH ₂) ₃ -NMe ₂	VIII.10 XV.268	C ₃₁ H ₃₃ FN ₄ O ₄	545 [M+H] ⁺	101	0.10 (B)
9.50		-N(Me)-(CO)- (CH ₂) ₂ -NMe ₂	VIII.10 XV.126	C ₃₀ H ₃₁ FN ₄ O ₄	531 [M+H] ⁺	161	0.20 (B)

9.51			VIII.17 XV.204	$C_{30}H_{31}FN_4O_4$	586 [M+H] ⁺	181- 183	0.20 (B)
9.52		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	VIII.17 XV.2	$C_{30}H_{33}FN_4O_5S$	581 [M+H] ⁺	158- 160	0.35 (B)
9.53		-N(Me)-(CO)- CH ₂ -NMe ₂	VIII.17 XV.125	$C_{30}H_{31}FN_4O_4$	531 [M+H] ⁺	n. d.	0.40 (B)
9.54		-N(COMe)- (CH ₂) ₃ -NMe ₂	VIII.17 XV.7	$C_{32}H_{35}FN_4O_4$	559 [M+H] ⁺	n. d.	0.50 (E)
9.55			VIII.11 XV.204	$C_{35}H_{41}FN_6O_4$	629 [M+H] ⁺	n. d.	0.35 (A)
9.56		-NMe-(CO)-CH ₃	VII.26 -	$C_{27}H_{25}FN_4O_3$	473 [M+H] ⁺	122- 126	0.50 (F)
9.57		-N(COMe)- (CH ₂) ₃ -NMe ₂	VII.26 XV.7	$C_{31}H_{34}FN_5O_3$	544 [M+H] ⁺	80-83	0.25 (A)
9.58		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	VII.26 XV.2	$C_{29}H_{32}FN_5O_4S$	566 [M+H] ⁺	190- 195	0.30 (A)

9.59		-N(Me)-(CO)- CH ₂ -NMe ₂	VII.26 XV.125	C ₂₉ H ₃₀ FN ₅ O ₃	516 [M+H] ⁺	238- 241	0.30 (G)
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*Eluant mixtures:

(A): silica gel, methylene chloride/methanol/ammonia 9:1:0.1

(B): silica gel, methylene chloride/methanol 9:1

(C): silica gel, methylene chloride/methanol/ammonia 8:1:0.1

(D): silica gel, methylene chloride/methanol/ammonia 10:1:0.1

(E): silica gel, methylene chloride/methanol/ammonia 5:1:0.01

(F): silica gel, ethyl acetate/methanol/ammonia = 9:1:0.1

(G): aluminium oxide, methylene chloride/methanol = 19:1

Example 10.0

3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-cyano-2-indolinone

130 mg of 1-acetyl-3-(1-methoxy-1-(3,4-dimethoxy-phenyl)-methylene)-6-cyano-2-indolinone (educt VIII.5) and 58 mg of 4-(dimethylaminomethyl)-aniline (educt XV.4) are dissolved in 5 ml of dimethylformamide and stirred for 2 hours at 80°C. After cooling the solvent is eliminated and the residue is purified through a silica gel column with methylene chloride/methanol 9:1 as eluant.

Yield: 21 mg (12% of theory),

R_f value: 0.35 (silica gel, methylene chloride/methanol = 9:1)

M.p. 265 °C

C₂₇H₂₆N₄O₃

Example 11.03-Z-[1-(4-(*N*-methyl-*N*-methylsulphonyl-amino)-anilino)-1-(3-(2-methoxycarbonyl-vinyl)-phenyl)-methylene]-6-chloro-2-indolinone

580 mg of 3-Z-[1-(4-(*N*-methyl-*N*-methylsulphonyl-amino)-anilino)-1-(3-iodo-phenyl)-methylene]-6-chloro-2-indolinone (educt 4.0) and 140 ml methyl acrylate are dissolved in 20 ml acetonitrile and 11 ml of dimethylformamide and 11 mg of palladium(II)-acetate, 2 ml of triethylamine and 30 mg of tri-ortho-tolyl-phosphine are added. The solution is stirred for 10 hours at 90°C under nitrogen as protective gas. After cooling it is filtered through Celite, the solvent is eliminated and the residue is purified through a silica gel column with methylene chloride/methanol 20:1 as eluant.

Yield: 450 mg (84% of theory),

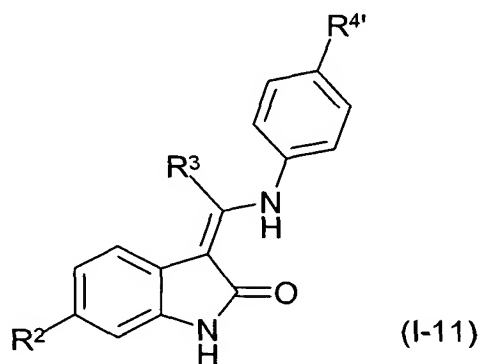
R_f value: 0.30 (silica gel, toluene/ethyl acetate = 1:1)

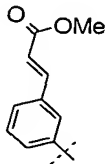
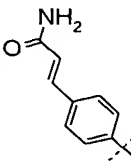
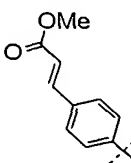
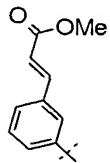
M.p. 228-232 °C

C₂₇H₂₄ClN₃O₅S

Mass spectrum: m/z = 537/539 [M]⁺

The following compounds of general formula I-11 are prepared analogously to Example 11.0:



Example	R ²	R ³	R ⁴	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
11.1	-Cl		-CH ₂ -NMe ₂	4.1	C ₂₈ H ₂₆ ClN ₃ O ₃	486/488 [M-H] ⁻	150-155	0.50 (A)
11.2	-F		-CH ₂ -NMe ₂	9.0	C ₂₇ H ₂₅ FN ₄ O ₂	455 [M-H] ⁻	269-270	0.20 (B)
11.3	-F		-CH ₂ -NMe ₂	9.0	C ₂₈ H ₂₆ FN ₃ O ₃	470 [M-H] ⁻	205-208	0.65 (A)
11.4	-F		-CH ₂ -NMe ₂	4.1	C ₂₈ H ₂₆ FN ₃ O ₃	472 [M+H] ⁺	138-140	0.45 (A)

*Eluant mixtures:

(A): silica gel, methylene chloride/methanol 5:1

(B): silica gel, methylene chloride/methanol/ammonia 9:1:0.01

Example 12.0

3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone

1.0 g of 3-Z-[1-(4-(dimethylaminomethyl)-anilino)-1-(3-(2-methoxycarbonyl-vinyl)-phenyl)-methylene]-6-chloro-2-indolinone (educt 11.1) are dissolved in 100 ml of methanol and 200 mg of 10% palladium/charcoal is added as catalyst. Then the mixture is hydrogenated for 6 hours at ambient temperature under 50 psi of hydrogen pressure. After the reaction has ended the catalyst

is filtered off, the solvent is eliminated and the residue is dried in vacuo at 100°C.

Yield: 900 mg (90% of theory),

R_f value: 0.40 (silica gel, methylene chloride/methanol = 9:1)

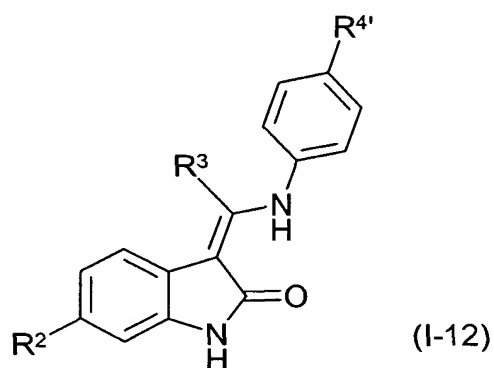
M.p. 160 °C

C₂₈H₂₈ClN₃O₃

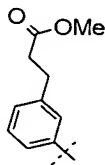
Mass spectrum: m/z = 490/492 [M+H]⁺

The following compounds of general formula I-12 are prepared analogously to

Example 12.0:



Example	R ²	R ³	R ⁴	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
12.1	-Cl		-N(Me)-SO ₂ Me	11.0	C ₂₇ H ₂₆ ClN ₃ O ₅ S	538/540 [M-H] ⁻	148-150	0.50 (A)
12.2	-F		-CH ₂ -NMe ₂	11.2	C ₂₇ H ₂₇ FN ₄ O ₂	459 [M+H] ⁺	150	0.70 (B)
12.3	-F		-CH ₂ -NMe ₂	11.3	C ₂₈ H ₂₈ FN ₃ O ₃	474 [M+H] ⁺	140	0.35 (A)

12.4	-F		-CH ₂ -NMe ₂	11.4	C ₂₈ H ₂₈ FN ₃ O ₃	474 [M+H] ⁺	140- 142	0.30 (A)
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*Eluant mixtures:

(A): silica gel, methylene chloride/methanol 9:1

(B): silica gel, methylene chloride/methanol/ammonia 5:1:0.01

Example 13.0

3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-amino-phenyl)-methylene]-6-chloro-2-indolinone

130 mg of 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-nitro-phenyl)-methylene]-6-chloro-2-indolinone (educt 5.18) are dissolved in 20 ml of ethanol and 20 ml of ethyl acetate and 100 mg of Raney nickel are added as catalyst. Then the mixture is hydrogenated for 20 hours at ambient temperature. After the end of the reaction the catalyst is filtered off, the solvent is eliminated, the residue is washed with a little diisopropylether and purified through a silica gel column with methylene chloride/ethanol/ammonia 30:1:0.1 as eluant. The product is washed with a little diisopropylether and dried in vacuo.

Yield: 80 mg (66% of theory),

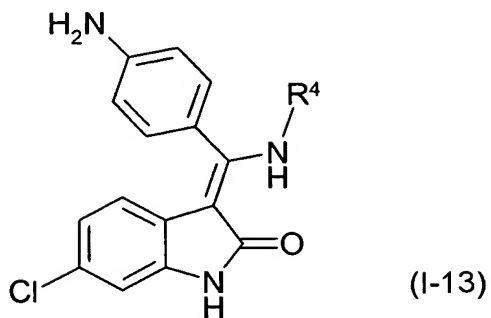
R_f value: 0.60 (silica gel, methylene chloride/ethanol/ammonia = 20:1:0.1)

M.p. 263-264 °C

C₂₄H₂₃ClN₄O

Mass spectrum: m/z = 419/421 [M+H]⁺

The following compounds of general formula I-13 are prepared analogously to Example 13.0 :



Example	R ⁴	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
13.1		-N(Me)-(CO)-CH ₂ -NMe ₂	5.21	C ₂₆ H ₂₆ ClN ₅ O ₂	476/478 [M+H] ⁺	275-276	0.10 (A)
13.2			5.20	C ₂₉ H ₃₁ ClN ₆ O ₂	529/531 [M-H] ⁻	268-269	0.15 (A)
13.3		-N(COMe)-(CH ₂) ₂ -NMe ₂	5.19	C ₂₇ H ₂₈ ClN ₅ O ₂	490/492 [M+H] ⁺	270	0.25 (A)
13.4			5.22	C ₂₇ H ₂₆ ClN ₅ O ₂	488/490 [M+H] ⁺	279-280	0.30 (A)

*Eluant mixtures:

(A): silica gel, methylene chloride/ethanol/ammonia 20:1:0.1

Example 14.03-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-aminomethyl-phenyl)-methylene]-6-chloro-2-indolinone

900 mg of 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-cyano-phenyl)-methylene]-6-chloro-2-indolinone (educt 5.23) are dissolved in 20 ml of methylene chloride, 30 ml of methanolic ammonia are added and 200 mg of Raney nickel are added as catalyst. Then the mixture is hydrogenated for 2 hours 15 minutes at ambient temperature and under 50 psi hydrogen pressure. After the end of the reaction the catalyst is filtered off, the solvent is eliminated and the residue is washed with a little methanol and diethyl ether. To liberate the base the residue is taken up in 1N sodium hydroxide solution and extracted four times with methylene chloride/methanol 9:1. The combined organic phases are washed with water and dried over sodium sulphate. The product is washed with a little diethyl ether and dried in vacuo.

Yield: 680 mg (75% of theory),

R_f value: 0.60 (silica gel, methylene chloride/methanol/ammonia = 9:1:0.1)

M.p. 211-214 °C

C₂₅H₂₅ClN₄O

Mass spectrum: m/z = 433/435 [M+H]⁺

Example 15.03-Z-[1-(4-(N-((4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-aminomethyl-phenyl)-methylene]-6-chloro-2-indolinone

1.39 g of 1-acetyl-3-Z-[1-(4-(N-((4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-cyano-phenyl)-methylene]-6-chloro-2-indolinone are dissolved in 20 ml methylene chloride, 30 ml of methanolic ammonia are added and 200 mg Raney nickel are added as catalyst. Then the mixture is hydrogenated for 2 hours at ambient temperature under 50 psi hydrogen pressure. After the end of the reaction the catalyst is filtered off, the solvent is eliminated and the residue is washed with a little methanol and diethyl ether. To liberate the base the residue is taken up in 1N sodium hydroxide solution and extracted four times with methylene chloride/methanol 9:1. The combined

organic phases are washed with water and dried over sodium sulphate. The product is purified through a silica gel column with a gradient of methylene chloride and methylene chloride/methanol/ammonia 8:1:0,1 as eluant. The product is washed with a little diethyl ether and dried in vacuo.

Yield: 700 mg (54% of theory),

R_f value: 0.15 (silica gel, methylene chloride/methanol/ammonia = 9:1:0.1)

M.p. 232-235 °C

C₃₀H₃₃ClN₆O₂

Mass spectrum: m/z = 544/546 [M]⁺

Example 16.0

3-Z-[1-(4-(N-((4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-hydroxy-phenyl)-methylene]-6-chloro-2-indolinone

100 mg of 3-Z-[1-(4-(N-((4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-benzyloxy-phenyl)-methylene]-6-chloro-2-indolinone (educt 4.9) are dissolved in 2 ml trifluoroacetic acid and stirred for 5 hours at 50°C. After the end of the reaction the solvent is eliminated. To liberate the base the residue is taken up in water and concentrated ammonia is added until an alkaline reaction is obtained. The precipitate formed is suction filtered, washed with water and at 100°C dried.

Yield: 10 mg (12% of theory),

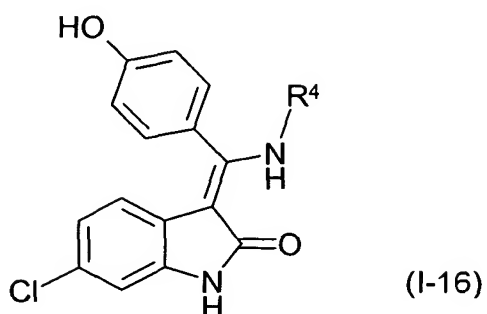
R_f value: 0.30 (silica gel, methylene chloride/methanol = 4:1)

M.p. 174-176 °C

C₂₉H₃₀ClN₅O₃

Mass spectrum: m/z = 532/534 [M+H]⁺

The following compounds of general formula I-16 are prepared analogously to Example 16.0:



Example	R ⁴	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
16.1		-N(Me)-(CO)-CH ₂ -NMe ₂	4.8	C ₂₆ H ₂₅ ClN ₄ O ₃	477/479 [M+H] ⁺	239-241	0.50 (A)
16.2		-N(COMe)-(CH ₂) ₂ -NMe ₂	4.10	C ₂₇ H ₂₇ ClN ₄ O ₃	491/493 [M+H] ⁺	249-251	0.40 (A)
16.3		-N(COMe)-(CH ₂) ₃ -NMe ₂	4.11	C ₂₈ H ₂₉ ClN ₄ O ₃	503/505 [M-H] ⁻	169-170	0.30 (A)
16.4		-CH ₂ -NMe ₂	4.7	C ₂₄ H ₂₂ ClN ₃ O ₂	418/420 [M-H] ⁻	215-217	0.05 (B)

*Eluant mixtures:

(A): silica gel, methylene chloride/methanol 4:1

(A): silica gel, methylene chloride/ethanol 10:1

Example 17.03-Z-[1-(4-ethylaminomethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone-trifluoroacetate

180 mg of 3-Z-[1-(4-(N-tert.butoxycarbonyl-ethylaminomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone (educt 1.30) are dissolved in 5 ml methylene chloride and 0.5 ml of trifluoroacetic acid are added. The mixture is stirred for 10 hours at ambient temperature. After this time the solvent is largely eliminated and the precipitate formed is suction filtered.

Yield: 110 mg (60% of theory),

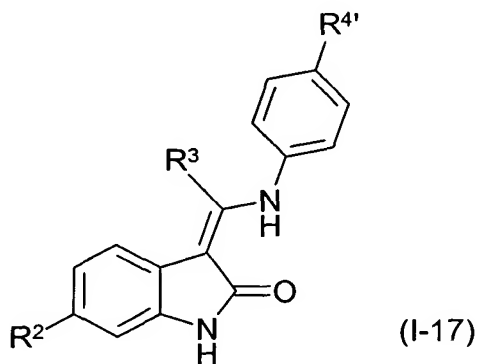
R_f value: 0.20 (silica gel, methylene chloride/methanol = 9:1)



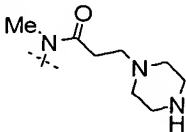

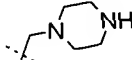

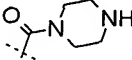


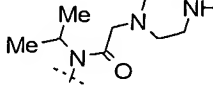

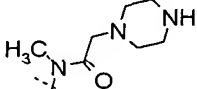

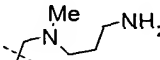
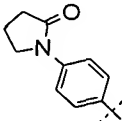


M.p. 260 °C



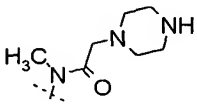


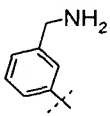
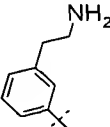
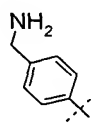
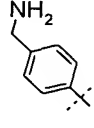
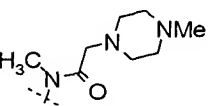
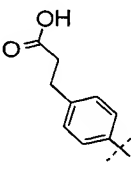
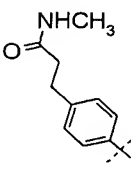
C₂₄H₂₂ClN₃O

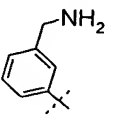
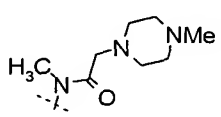
Mass spectrum: m/z = 402/404 [M-H]⁻

The following compounds of general formula I-17 are prepared analogously to Example 17.0 :



Example	R ²	R ³	R ⁴	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
17.1	-Cl		-CH ₂ -NHMe	1.43	C ₂₃ H ₂₀ ClN ₃ O	388/390 [M-H] ⁻	250	0.15 (A)
17.2	-Cl			1.50	C ₂₉ H ₃₀ ClN ₅ O ₂	514/516 [M-H] ⁻	224	0.25 (A)
17.3	-Cl			1.52	C ₂₆ H ₂₅ ClN ₄ O	443/445 [M-H] ⁻	240	0.25 (A)
17.4	-Cl			1.67	C ₂₆ H ₂₃ ClN ₄ O ₂	457/459 [M-H] ⁻	289	0.25 (A)
17.5	-Cl		-CH ₂ -NH ₂	1.77	C ₂₂ H ₁₈ ClN ₃ O	374/376 [M-H] ⁻	265	0.70 (B)
17.6	-Cl			1.79	C ₃₀ H ₃₂ ClN ₅ O ₂	528/530 [M-H] ⁻	164	0.70 (B)
17.7	-Cl			1.84	C ₂₈ H ₂₈ ClN ₅ O ₂	500/502 [M-H] ⁻	172	0.70 (B)
17.8	-Cl			1.87	C ₂₆ H ₂₇ ClN ₄ O	447/449 [M+H] ⁺	221	0.70 (B)
17.9	-Cl		-CH ₂ -NHMe	5.11	C ₂₇ H ₂₅ ClN ₄ O ₂	473/475 [M+H] ⁺	240- 244	0.25 (C)
17.10	-Br		-CH ₂ -NHEt	6.5	C ₂₄ H ₂₂ BrN ₃ O	446/448 [M-H] ⁻	274- 276	0.10 (A)
17.11	-Br		-CH ₂ -NHMe	6.17	C ₂₃ H ₂₀ BrN ₃ O	456/458 [M+Na] ⁺	252- 255	0.40 (C)

17.12	-CN		-CH ₂ -NHEt	7.5	C ₂₅ H ₂₂ N ₄ O	393 [M-H] ⁻	249	0.25 (A)
17.13	-CN			7.18	C ₂₉ H ₂₈ N ₆ O ₂	491 [M-H] ⁻	146	0.15 (A)
17.14	-F		-CH ₂ -NHEt	8.4	C ₂₄ H ₂₂ FN ₃ O	386 [M-H] ⁻	285- 288	0.40 (D)
17.15	-F		-CH ₂ -NHMe	8.16	C ₂₃ H ₂₀ FN ₃ O	372 [M-H] ⁻	251	0.15 (C)
17.16	-F		-CH ₂ -NMe ₂	9.13	C ₂₅ H ₂₅ FN ₄ O	415 [M-H] ⁻	168- 175	0.25 (C)
17.17	-F		-CH ₂ -NMe ₂	9.19	C ₂₆ H ₂₇ FN ₄ O	431 [M+H] ⁺	155- 160	0.45 (E)
17.18	-F		-CH ₂ -NMe ₂	9.18	C ₂₅ H ₂₅ FN ₄ O	417 [M+H] ⁺	203- 207	0.25 (C)
17.19	-F			9.17	C ₃₀ H ₃₃ FN ₆ O ₂	529 [M+H] ⁺	170- 175	0.15 (C)
17.20	-F		-CH ₂ -NHMe	20.11	C ₂₆ H ₂₄ FN ₃ O ₃	446 [M+H] ⁺	245- 251	0.20 (F)
17.21	-F		-CH ₂ -NHMe	21.22	C ₂₆ H ₂₄ FN ₃ O ₃	459 [M+H] ⁺	239- 243	0.30 (C)

17.22	-F			9.55	$C_{30}H_{33}FN_6O_2$	529 [M+H] ⁺	n. d.	n. d.
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*Eluant mixtures:

(A): silica gel, methylene chloride/methanol 9:1

(B): silica gel, methylene chloride/methanol 10:1

(C): silica gel, methylene chloride/methanol/ammonia 9:1:0.1

(D): silica gel, methylene chloride/methanol/ammonia 9:1:0.01

(E): silica gel, methylene chloride/methanol/ammonia 8:2:0.2

(F): Reversed phase RP8, methanol/saline solution (5%) = 3:2

Example 18.0

3-Z-[1-(4-(N-aminomethylcarbonyl)-N-methyl-amino)-anilino]-1-phenyl-methylene]-6-chloro-2-indolinone

600 mg of 3-Z-[1-(4-(N-(phthalimido-2-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone (educt 1.42) are dissolved in 30 ml of ethanol and 500 mg of hydrazine hydrate (80%) are added. The mixture is stirred for 4 hours at 50°C. After cooling 20 ml methylene chloride are added and the precipitate formed is suction filtered. The filtrate is concentrated by evaporation and purified through a silica gel column with methylene chloride/methanol/ammonia 10:1:0.1 as eluant.

Yield: 100 mg (22% of theory),

R_f value: 0.40 (silica gel, methylene chloride/methanol/ammonia = 9:1:0.1)

M.p. 204-206 °C

$C_{24}H_{21}ClN_4O_2$

Mass spectrum: m/z = 432/434 [M]⁺

Example 19.03-Z-[1-(4-carboxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

450 mg of 3-Z-[1-(4-ethoxycarbonyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone (educt 1.2) are dissolved in 10 ml of ethanol and 2 ml of 1N sodium hydroxide solution are added. The mixture is stirred for 3 hours at 80°C. After cooling, 2 ml of 1N hydrochloric acid are added and the mixture is stirred for half an hour at ambient temperature. The precipitate formed is suction filtered and washed with ethanol and diethyl ether.

Yield: 320 mg (76% of theory),

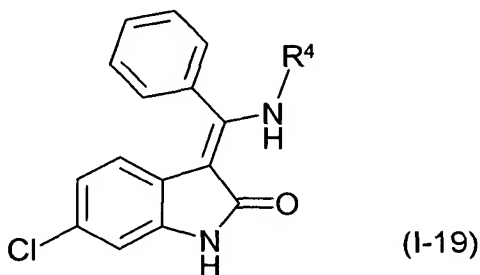
R_f value: 0.40 (silica gel, methylene chloride/ethanol = 9:1)

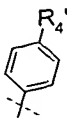
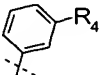
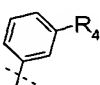
M.p. 333-334 °C (decomposition)

C₂₂H₁₅ClN₂O₃

Mass spectrum: m/z = 389/391 [M-H]⁻

The following compounds of general formula I-19 are prepared analogously to Example 19.0 :



Example	R ⁴	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
19.1		-CH ₂ -COOH	1.100	C ₂₃ H ₁₇ ClN ₂ O ₃	403/405 [M-H] ⁻	277	0.25 (A)
19.2		-CH ₂ -COOH	1.74	C ₂₃ H ₁₇ ClN ₂ O ₃	403/405 [M-H] ⁻	209	0.15 (A)
19.3		-COOH	1.75	C ₂₂ H ₁₅ ClN ₂ O ₃	389/391 [M-H] ⁻	321	0.25 (A)

*Eluant mixtures:

(A): silica gel, methylene chloride/methanol 9:1

Example 20.0

3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone

900 mg of 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-methoxycarbonyl-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone (educt 12.0) are dissolved in 10 ml of ethanol and 5 ml of 1N sodium hydroxide solution are added. The mixture is stirred for 5 hours at ambient temperature. After cooling, 5 ml of 1N hydrochloric acid are added. The precipitate formed is suction filtered and washed with water.

Yield: 830 mg (95% of theory),

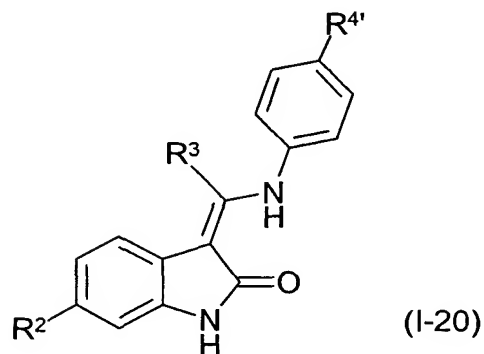
R_f value: 0.50 (Reversed phase RP8, methanol/saline solution (5%) = 4:1)

M.p. 210-215 °C

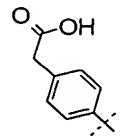
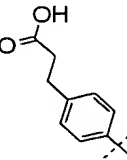
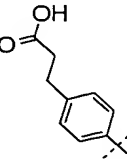
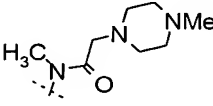
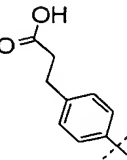
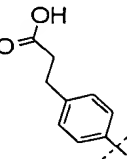
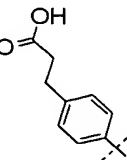
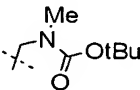
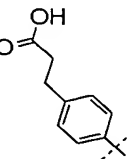
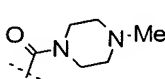
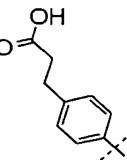

C₂₇H₂₆ClN₃O₃

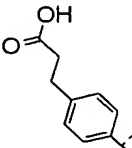
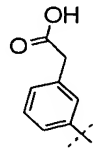
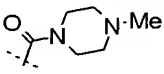
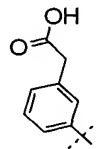
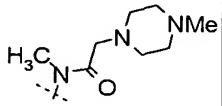
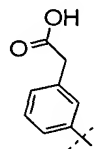
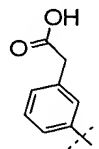
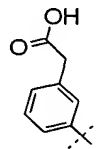
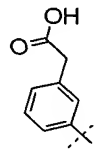
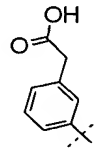

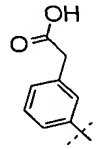
Mass spectrum: m/z = 476/478 [M+H]⁺

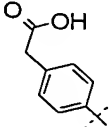
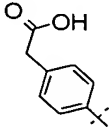

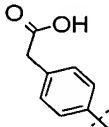
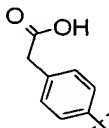
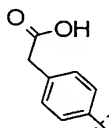
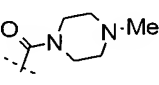
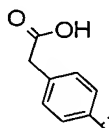
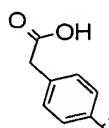
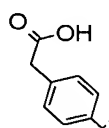
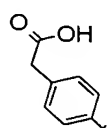
The following compounds of general formula I-20 are prepared analogously to Example 20.0 :

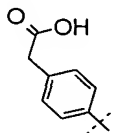
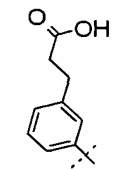
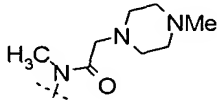
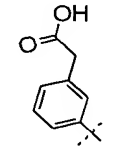
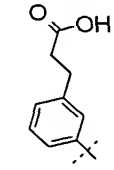
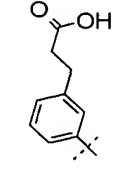
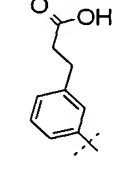


Example	R ²	R ³	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
20.1	-F		-CH ₂ -NMe ₂	12.3	C ₂₇ H ₂₆ FN ₃ O ₃	460 [M+H] ⁺	250	0.65 (A)
20.2	-F		-CH ₂ -NMe ₂	9.12	C ₂₆ H ₂₄ FN ₃ O ₃	444 [M-H] ⁻	278- 282	0.10 (B)
20.3	-F		-CH ₂ -NMe ₂	12.4	C ₂₇ H ₂₆ FN ₃ O ₃	458 [M-H] ⁻	198- 200	0.20 (C)
20.4	-F		-CH ₂ -NMe ₂	9.10	C ₂₆ H ₂₄ FN ₃ O ₃	444 [M-H] ⁻	212- 216	0.30 (D)
20.5	-F			9.15	C ₃₁ H ₃₂ FN ₅ O ₄	558 [M+H] ⁺	260- 263	0.20 (D)

20.6	-F		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	9.14	C ₂₈ H ₂₉ FN ₄ O ₅ S	553 [M+H] ⁺	246- 249	0.30 (D)
20.7	-F		-NMe-(CO)- CH ₃	9.20	C ₂₇ H ₂₄ FN ₃ O ₄	474 [M+H] ⁺	286- 290	0.60 (E)
20.8	-F			9.21	C ₃₂ H ₃₄ FN ₅ O ₄	570 [M-H] ⁻	215- 222	0.20 (D)
20.9	-F		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	9.22	C ₂₉ H ₃₁ FN ₄ O ₅ S	567 [M+H] ⁺	160- 165	0.20 (D)
20.10	-F		-N(COMe)- (CH ₂) ₃ -NMe ₂	9.23	C ₃₁ H ₃₃ FN ₄ O ₄	545 [M+H] ⁺	153- 158	0.15 (D)
20.11	-F			9.24	C ₃₁ H ₃₂ FN ₃ O ₅	546 [M+H] ⁺	215- 219	0.60 (E)
20.12	-F			9.25	C ₃₀ H ₂₉ FN ₄ O ₄	529 [M+H] ⁺	179- 186	0.25 (E)
20.13	-F			9.26	C ₂₈ H ₂₃ FN ₄ O ₃	483 [M+H] ⁺	264- 267	0.65 (E)

20.14	-F		-SO ₂ Me	9.27	C ₂₅ H ₂₁ FN ₂ O ₅ S	481 [M+H] ⁺	146- 155	0.70 (E)
20.15	-F			9.30	C ₂₉ H ₂₇ FN ₄ O ₄	515 [M+H] ⁺	251	0.70 (E)
20.16	-F			9.28	C ₃₁ H ₃₂ FN ₅ O ₄	558 [M+H] ⁺	234	0.10 (E)
20.17	-F		-N(Me)-(CO)- CH ₂ -NMe ₂	9.31	C ₂₈ H ₂₇ FN ₄ O ₄	503 [M+H] ⁺	203	0.60 (E)
20.18	-F		-N(Me)-(CO)- (CH ₂) ₄ -NMe ₂	9.34	C ₃₁ H ₃₃ FN ₄ O ₄	545 [M+H] ⁺	251	n. d.
20.19	-F		-H	9.45	C ₂₃ H ₁₇ FN ₂ O ₃	387 [M-H] ⁻	130	0.60 (E)
20.20	-F		-SO ₂ Me	9.46	C ₂₄ H ₁₉ FN ₂ O ₅ S	467 [M+H] ⁺	139	0.55 (E)
20.21	-F			9.47	C ₂₇ H ₂₁ FN ₄ O ₃	469 [M+H] ⁺	157	0.35 (E)
20.22	-F		-N(SO ₂ Me)- (CH ₂)-(CO)- NMe ₂	9.48	C ₂₈ H ₂₇ FN ₄ O ₆ S	567 [M+H] ⁺	183	0.55 (E)

20.23	-F		-H	9.35	$C_{23}H_{17}FN_2O_3$	389 [M+H] ⁺	237- 240	0.10 (D)
20.24	-F			9.36	$C_{27}H_{21}FN_4O_3$	469 [M+H] ⁺	259- 265	0.15 (D)
20.25	-F		-N(COMe)- (CH ₂) ₃ -NMe ₂	9.44	$C_{30}H_{31}FN_4O_4$	531 [M+H] ⁺	274- 278	0.15 (D)
20.26	-F		-N(Me)-(CO)- CH ₂ -NMe ₂	9.39	$C_{28}H_{27}FN_4O_4$	503 [M+H] ⁺	258- 264	0.20 (D)
20.27	-F			9.37	$C_{29}H_{27}FN_4O_4$	515 [M+H] ⁺	279- 282	0.15 (D)
20.28	-F		-SO ₂ Me	9.42	$C_{24}H_{19}FN_2O_5S$	467 [M+H] ⁺	260- 266	0.35 (F)
20.29	-F		-N(COMe)- CH ₃	9.40	$C_{26}H_{22}FN_3O_4$	460 [M+H] ⁺	290- 294	0.30 (F)
20.30	-F		-N(SO ₂ Me)- CH ₂ -(CO)- NMe ₂	9.38	$C_{28}H_{27}FN_4O_6S$	567 [M+H] ⁺	238- 242	0.30 (F)
20.31	-F		-N(Me)-(CO)- (CH ₂) ₂ -NMe ₂	9.41	$C_{29}H_{29}FN_4O_4$	517 [M+H] ⁺	250- 255	0.35 (F)

20.32	-F		-N(Me)-(CO)- (CH ₂) ₃ -NMe ₂	9.43	C ₃₀ H ₃₁ FN ₄ O ₄	531 [M+H] ⁺	184- 190	0.25 (F)
20.33	-F			9.51	C ₃₂ H ₃₄ FN ₅ O ₄	572 [M-H] ⁻	170- 175	0.40 (C)
20.34	-F		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	9.29	C ₂₈ H ₂₉ FN ₄ O ₅ S	553 [M+H] ⁺	180	0.60 (C)
20.35	-F		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	9.52	C ₂₉ H ₃₁ FN ₄ O ₅ S	567 [M+H] ⁺	196- 199	0.30 (C)
20.36	-F		-N(Me)-(CO)- CH ₂ -NMe ₂	9.53	C ₂₉ H ₂₉ FN ₄ O ₄	517 [M+H] ⁺	150	0.20 (C)
20.37	-F		-N(COMe)- (CH ₂) ₃ -NMe ₂	9.54	C ₃₁ H ₃₃ FN ₄ O ₄	545 [M+H] ⁺	206- 210	0.30 (A)

*Eluant mixtures:

(A): Reversed phase RP8, methanol/saline solution(5%) = 4:1

(B): silica gel, methylene chloride/methanol = 8:2

(C): silica gel, methylene chloride/methanol = 5:1

(D): Reversed phase RP8, methanol/saline solution(5%) = 3:2

(E): silica gel, methylene chloride/methanol = 9:1

(F): Reversed phase RP8, methanol/saline solution(5%) = 7:3

Example 21.03-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carbamoyl-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone

480 mg of 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(3-(2-carboxyethyl)-phenyl)-methylene]-6-chloro-2-indolinone (educt 20.0), 350 mg TBTU, 150 mg HOBT and 420 ml triethylamine are dissolved in 10 ml of dimethylformamide and 620 mg of N-hydroxysuccinimide-ammonium salt are added. The mixture is stirred for 20 hours at ambient temperature. After the solvent has been eliminated the residue is suspended in a little ethyl acetate and water, filtered off and washed with water. The residue is purified through an aluminium oxide column (activity 2-3) with methylene chloride/ethanol 20:1 as eluant. The product is recrystallised from diethyl ether and dried in vacuo at 100°C.

Yield: 370 mg (78% of theory),

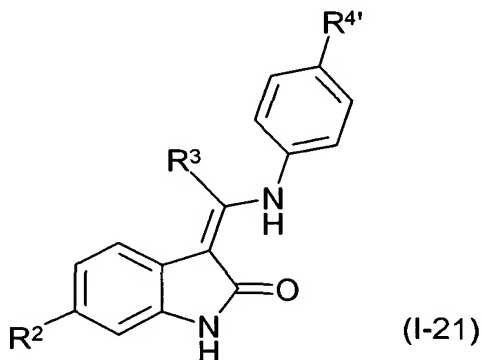
R_f value: 0.40 (aluminium oxide, methylene chloride/ethanol = 20:1)

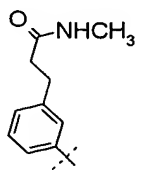
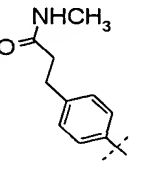
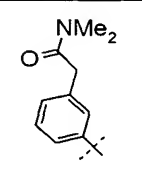
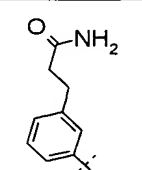
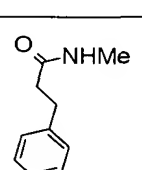
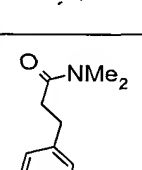
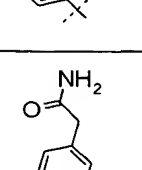
M.p. 222-225 °C

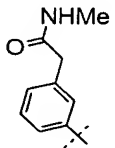
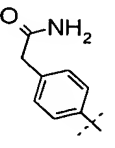
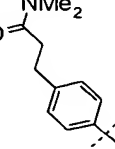
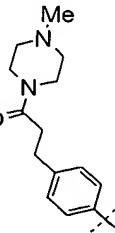
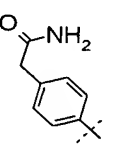
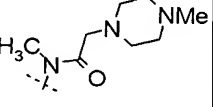
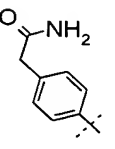
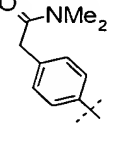
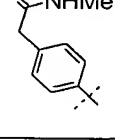
C₂₇H₂₇ClN₄O₂

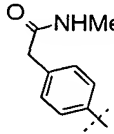
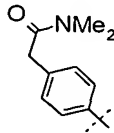
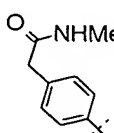
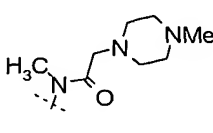
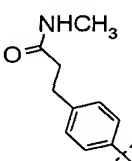
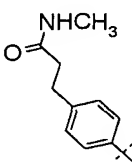
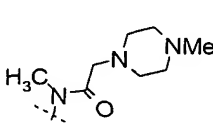
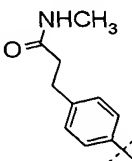
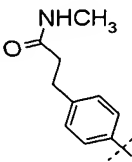
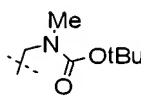
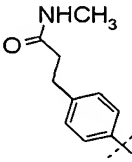

Mass spectrum: m/z = 475/477 [M+H]⁺

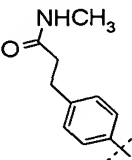
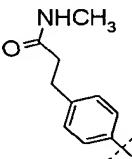
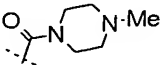
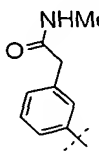
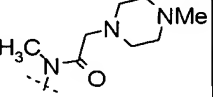
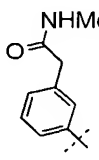
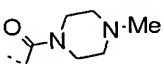
The following compounds of general formula I-21 are prepared analogously to Example 21.0 :



Example	R ²	R ³	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
21.1	-Cl		-CH ₂ -NMe ₂	20.0 **	C ₂₈ H ₂₉ ClN ₄ O ₂	489/491 [M+H] ⁺	223- 225	0.50 (A)
21.2	-F		-CH ₂ -NMe ₂	20.1 **	C ₂₈ H ₂₉ FN ₄ O ₂	473 [M+H] ⁺	148- 150	0.40 (B)
21.3	-F		-CH ₂ -NMe ₂	20.2 ***	C ₂₈ H ₂₉ FN ₄ O ₂	473 [M+H] ⁺	98- 103	0.30 (C)
21.4	-F		-CH ₂ -NMe ₂	20.3	C ₂₇ H ₂₇ FN ₄ O ₂	459 [M+H] ⁺	223- 225	0.50 (A)
21.5	-F		-CH ₂ -NMe ₂	20.3 **	C ₂₈ H ₂₉ FN ₄ O ₂	473 [M+H] ⁺	210- 213	0.70 (A)
21.6	-F		-CH ₂ -NMe ₂	20.3 ***	C ₂₉ H ₃₁ FN ₄ O ₂	487 [M+H] ⁺	213- 215	0.80 (A)
21.7	-F		-CH ₂ -NMe ₂	20.2	C ₂₆ H ₂₅ FN ₄ O ₂	443 [M-H] ⁻	115- 120	0.25 (C)

21.8	-F		-CH ₂ -NMe ₂	20.2 **	C ₂₇ H ₂₇ FN ₄ O ₂	457 [M-H] ⁻	222- 225	0.25 (C)
21.9	-F		-CH ₂ -NMe ₂	20.4	C ₂₆ H ₂₅ FN ₄ O ₂	443 [M-H] ⁻	143- 146	0.40 (D)
21.10	-F		-CH ₂ -NMe ₂	20.1 ***	C ₂₉ H ₃₁ FN ₄ O ₂	487 [M+H] ⁺	198- 200	0.60 (B)
21.11	-F		-CH ₂ -NMe ₂	20.1 ****	C ₃₂ H ₃₆ FN ₅ O ₂	542 [M+H] ⁺	175	0.60 (B)
21.12	-F			20.5	C ₃₁ H ₃₃ FN ₆ O ₃	557 [M+H] ⁺	150- 156	0.40 (E)
21.13	-F		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	20.6	C ₂₈ H ₃₀ FN ₅ O ₄ S	552 [M+H] ⁺	197- 199	0.50 (D)
21.14	-F		-CH ₂ -NMe ₂	20.4 ***	C ₂₈ H ₂₉ FN ₄ O ₂	473 [M+H] ⁺	147- 152	0.35 (D)
21.15	-F		-CH ₂ -NMe ₂	20.4 **	C ₂₇ H ₂₇ FN ₄ O ₂	459 [M+H] ⁺	208- 214	0.35 (D)

21.16	-F		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	20.6 **	C ₂₉ H ₃₂ FN ₅ O ₄ S	566 [M+H] ⁺	218- 222	0.70 (F)
21.17	-F		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	20.6 ***	C ₃₀ H ₃₄ FN ₅ O ₄ S	580 [M+H] ⁺	199- 205	0.40 (C)
21.18	-F			20.5 **	C ₃₂ H ₃₅ FN ₆ O ₃	571 [M+H] ⁺	155- 160	0.20 (C)
21.19	-F		-N(Me)-(CO)- CH ₃	20.7 **	C ₂₈ H ₂₇ FN ₄ O ₃	487 [M+H] ⁺	137- 145	0.50 (C)
21.20	-F			20.8 **	C ₃₃ H ₃₇ FN ₆ O ₃	585 [M+H] ⁺	211- 219	0.40 (C)
21.21	-F		-N(SO ₂ Me)- (CH ₂) ₂ -NMe ₂	20.9 **	C ₃₀ H ₃₄ FN ₅ O ₄ S	578 [M-H] ⁻	192- 200	0.50 (C)
21.22	-F			20.11 **	C ₃₂ H ₃₅ FN ₄ O ₄	559 [M+H] ⁺	180- 187	0.50 (C)
21.23	-F			20.13 **	C ₂₉ H ₂₆ FN ₅ O ₂	496 [M+H] ⁺	262- 266	0.40 (C)

21.24	-F		-SO ₂ Me	20.14 **	C ₂₆ H ₂₄ FN ₃ O ₄ S	494 [M+H] ⁺	180- 188	0.60 (C)
21.25	-F			20.12 **	C ₃₁ H ₃₂ FN ₅ O ₃	542 [M+H] ⁺	226- 230	0.50 (C)
21.26	-F			20.16 **	C ₃₂ H ₃₅ FN ₆ O ₃	571 [M+H] ⁺	213	0.10 (G)
21.27	-F			20.15 **	C ₃₀ H ₃₀ FN ₅ O ₃	528 [M+H] ⁺	245	0.40 (G)

*Eluant mixtures:

(A): silica gel, methylene chloride/methanol/ammonia = 5:1:0.01

(B): aluminium oxide, methylene chloride/ethanol = 20:1

(C): silica gel, methylene chloride/methanol/ammonia = 9:1:0.1

(D): silica gel, methylene chloride/methanol/ammonia = 6:1:0.1

(E): silica gel, methylene chloride/methanol/ammonia = 5:1:0.1

(F): silica gel, methylene chloride/methanol/ammonia = 7:1:0.1

(G): silica gel, methylene chloride/methanol = 9:1

** with methylammonium chloride as base equivalent

*** with dimethylammonium chloride as base equivalent

**** with piperidine hydrochloride as base equivalent

Example 22.03-Z-[1-(4-(4-methyl-piperazin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-1,3-dihydro-indol-2-thione

460 mg of 3-Z-[1-(4-(4-methyl-piperazin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone (educt 1.35) are dissolved in 5 ml of pyridine and 220 mg of phosphorus pentasulphide are added. The mixture is stirred for 2 hours at 120°C. After cooling it is diluted with water and 0.5 ml of concentrated ammonia is added. The precipitate formed is suction filtered and purified through a silica gel column with methylene chloride/methanol 9:1 as eluant. The product is recrystallised from petroleum ether and dried in vacuo at 100°C.

Yield: 300 mg (63% of theory),

R_f value: 0.50 (silica gel, methylene chloride/methanol = 9:1)

M.p. 250-252 °C

C₂₇H₂₇ClN₄S

Mass spectrum: m/z = 475/477 [M+H]⁺

Example 23.03-Z-[1-(4-(N-acetylaminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

61 mg of 3-Z-[1-(4-(N-aminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone (educt 18.0) are dissolved in 3 ml acetic acid and 0.1 ml acetic anhydride are added. The mixture is stirred for 1.5 hours at ambient temperature. After this time the solvent is eliminated, the residue is suspended in a little water and suction filtered. The product is dried in vacuo at 100°C.

Yield: 60 mg (90% of theory),

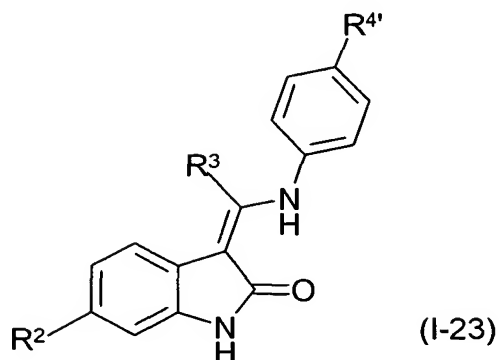
R_f value: 0.50 (silica gel, methylene chloride/methanol = 9:1)

M.p. 291-292 °C

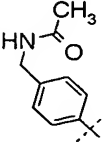
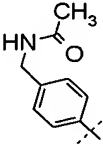
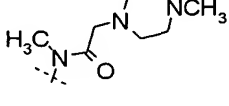
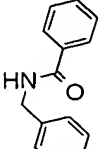
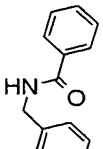
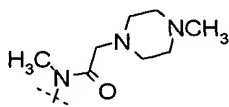
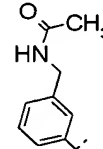
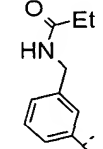
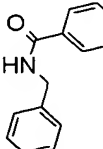
C₂₆H₂₃ClN₄O₃

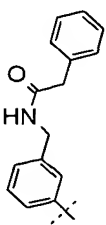
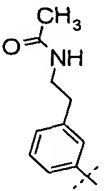
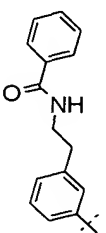
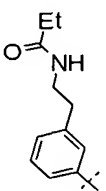
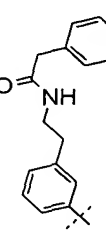
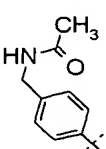
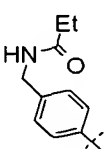
Mass spectrum: m/z = 497/499 [M+Na]⁺

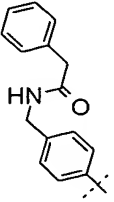
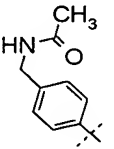
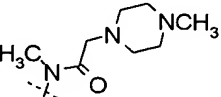
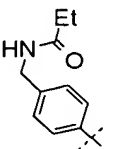
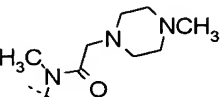
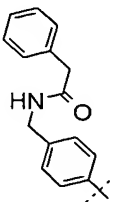
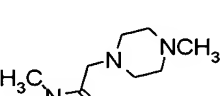
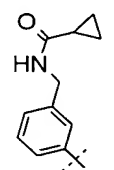
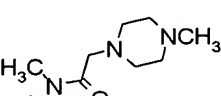
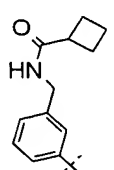
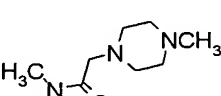
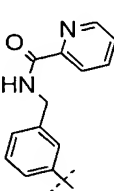
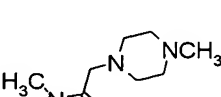
The following compounds of general formula I-23 are prepared analogously to Example 23.0:

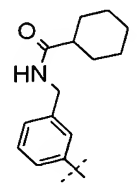
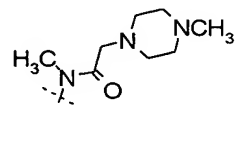
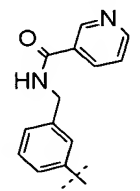
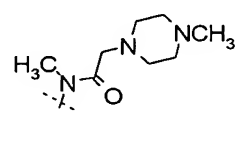
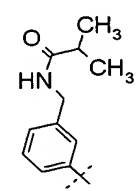
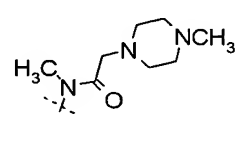
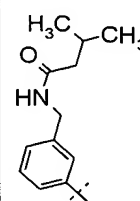
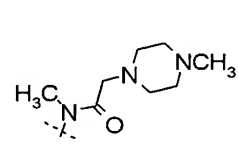
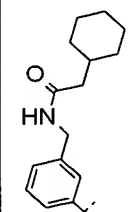
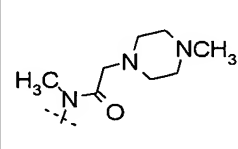
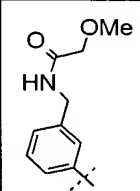
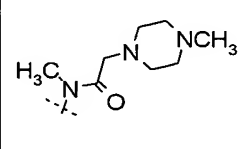
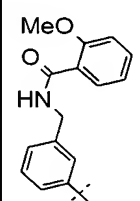
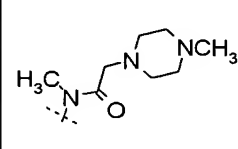


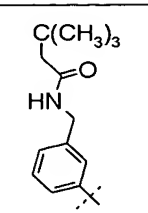
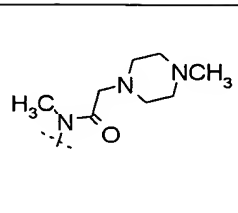
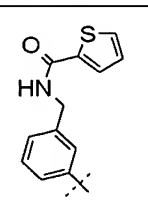
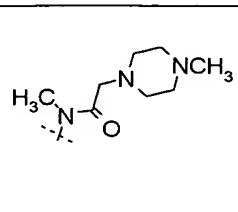
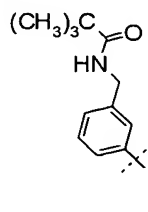
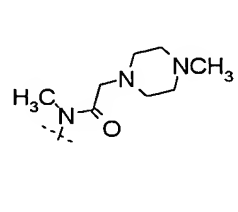
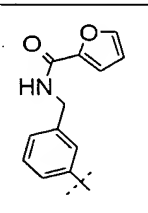
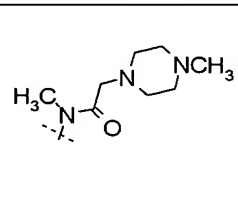
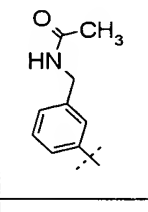
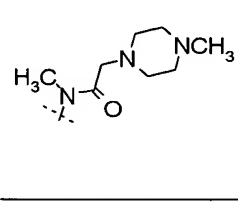
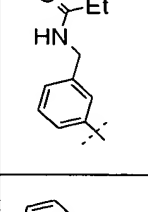
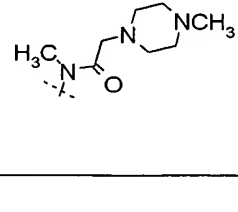
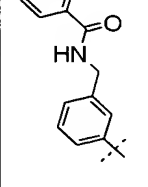
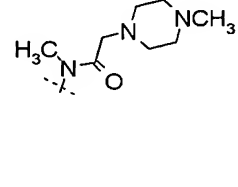
Example	R ²	R ³	R ⁴	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
23.1	-Cl		-N(COMe)-(CH ₂) ₂ -NMe ₂	13.3	C ₂₉ H ₃₀ ClN ₅ O ₃	530/532 [M-H] ⁻	187-188	0.15 (A)
23.2	-Cl		-N(Me)-(CO)-CH ₂ -NMe ₂	13.1	C ₂₈ H ₂₈ ClN ₅ O ₃	518/520 [M+H] ⁺	249-250	0.15 (A)
23.3	-Cl			13.2	C ₃₁ H ₃₃ ClN ₆ O ₃	571/573 [M-H] ⁻	168-170	0.10 (A)
23.4	-Cl			13.4	C ₂₉ H ₂₈ ClN ₅ O ₃	528/530 [M-H] ⁻	160	0.15 (A)
23.5	-Cl		-CH ₂ -NMe ₂	13.0	C ₂₆ H ₂₅ ClN ₄ O ₂	459/461 [M-H] ⁻	158-159	0.25 (A)

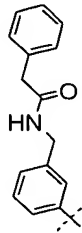
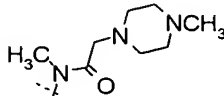
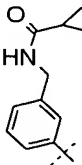
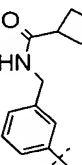
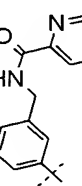
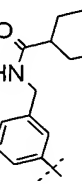
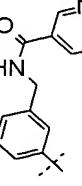
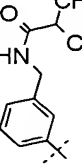
23.6	-Cl		-CH ₂ -NMe ₂	14.0	C ₂₇ H ₂₇ ClN ₄ O ₂	473/475 [M-H] ⁻	219- 220	0.30 (B)
23.7	-Cl			15.0	C ₃₂ H ₃₅ ClN ₆ O ₃	585/587 [M-H] ⁻	252- 255	0.25 (B)
23.8	-Cl		-CH ₂ -NMe ₂	14.0	C ₃₂ H ₂₉ ClN ₄ O ₂	535/537 [M-H] ⁻	238 (dec.)	0.45 (B)
23.9	-Cl			15.0	C ₃₇ H ₃₇ ClN ₆ O ₃	647/649 [M-H] ⁻	282- 284	0.40 (B)
23.10	-F		-CH ₂ -NMe ₂	17.16	C ₂₇ H ₂₇ FN ₄ O ₂	457 [M-H] ⁻	245- 250	0.40 (C)
23.11	-F		-CH ₂ -NMe ₂	17.16	C ₂₈ H ₂₉ FN ₄ O ₂	471 [M-H] ⁻	212- 214	0.35 (D)
23.12	-F		-CH ₂ -NMe ₂	17.16	C ₃₂ H ₂₉ FN ₄ O ₂	519 [M-H] ⁻	237- 240	0.40 (D)

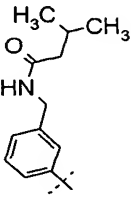
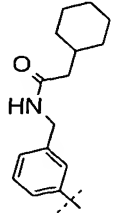
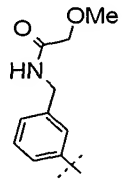
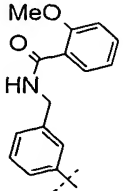
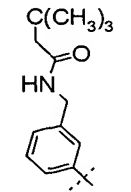
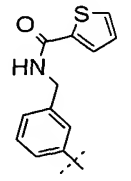
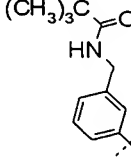
23.13	-F		-CH ₂ -NMe ₂	17.16	C ₃₃ H ₃₁ FN ₄ O ₂	533 [M-H] ⁻	187- 190	0.30 (D)
23.14	-F		-CH ₂ -NMe ₂	17.17	C ₂₈ H ₂₉ FN ₄ O ₂	471 [M-H] ⁻	234- 237	0.30 (D)
23.15	-F		-CH ₂ -NMe ₂	17.17	C ₃₃ H ₃₁ FN ₄ O ₂	533 [M-H] ⁻	144- 150	0.45 (C)
23.16	-F		-CH ₂ -NMe ₂	17.17	C ₂₉ H ₃₁ FN ₄ O ₂	485 [M-H] ⁻	235- 237	0.25 (D)
23.17	-F		-CH ₂ -NMe ₂	17.17	C ₃₄ H ₃₃ FN ₄ O ₂	547 [M-H] ⁻	217- 220	0.30 (D)
23.18	-F		-CH ₂ -NMe ₂	17.18	C ₂₇ H ₂₇ FN ₄ O ₂	457 [M-H] ⁻	112- 120	0.25 (D)
23.19	-F		-CH ₂ -NMe ₂	17.18	C ₂₈ H ₂₉ FN ₄ O ₂	586 [M+H] ⁺	176- 180	0.30 (D)

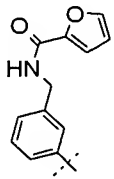
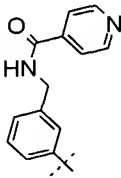
23.20	-F		-CH ₂ -NMe ₂	17.18	C ₃₃ H ₃₁ FN ₄ O ₂	535 [M+H] ⁺	80- 85	0.35 (D)
23.21	-F			17.19	C ₃₂ H ₃₅ FN ₆ O ₃	569 [M-H] ⁻	230- 235	0.35 (D)
23.22	-F			17.19	C ₃₃ H ₃₇ FN ₆ O ₃	583 [M-H] ⁻	205- 210	0.30 (D)
23.23	-F			17.19	C ₃₈ H ₃₉ FN ₆ O ₃	645 [M-H] ⁻	217- 220	0.35 (D)
23.24	-F			17.22	C ₃₄ H ₃₇ FN ₆ O ₃	597 [M+H] ⁺	209- 212	0.30 (D)
23.25	-F			17.22	C ₃₅ H ₃₉ FN ₆ O ₃	611 [M+H] ⁺	190- 193	0.30 (D)
23.26	-F			17.22	C ₃₆ H ₃₆ FN ₇ O ₃	634 [M+H] ⁺	160- 163	0.30 (D)

23.27	-F			17.22	$C_{37}H_{43}FN_6O_3$	639 [M+H] ⁺	223- 227	0.30 (D)
23.28	-F			17.22	$C_{36}H_{36}FN_7O_3$	634 [M+H] ⁺	170- 175	0.25 (D)
23.29	-F			17.22	$C_{34}H_{39}FN_6O_3$	599 [M+H] ⁺	194- 196	0.20 (D)
23.30	-F			17.22	$C_{35}H_{41}FN_6O_3$	613 [M+H] ⁺	197- 200	0.70 (E)
23.31	-F			17.22	$C_{38}H_{45}FN_6O_3$	653 [M+H] ⁺	130- 135	0.75 (E)
23.32	-F			17.22	$C_{33}H_{37}FN_6O_4$	601 [M+H] ⁺	155- 159	0.60 (E)
23.33	-F			17.22	$C_{38}H_{39}FN_6O_4$	663 [M+H] ⁺	168- 172	0.35 (C)

23.34	-F			17.22	$C_{36}H_{43}FN_6O_3$	627 [M+H] ⁺	85-90	0.35 (C)
23.35	-F			17.22	$C_{35}H_{35}FN_6O_3S$	639 [M+H] ⁺	170-175	0.25 (C)
23.36	-F			17.22	$C_{35}H_{41}FN_6O_3$	613 [M+H] ⁺	242-245	0.30 (C)
23.37	-F			17.22	$C_{35}H_{35}FN_6O_4$	623 [M+H] ⁺	155-160	0.65 (F)
23.38	-F			17.22	$C_{32}H_{35}FN_6O_3$	571 [M+H] ⁺	190-195	0.60 (F)
23.39	-F			17.22	$C_{33}H_{37}FN_6O_3$	585 [M+H] ⁺	203-209	0.65 (E)
23.40	-F			17.22	$C_{37}H_{37}FN_6O_3$	633 [M+H] ⁺	145-150	0.60 (F)

23.41	-F			17.22	$C_{38}H_{39}FN_6O_3$	647 [M+H] ⁺	148- 151	0.65 (F)
23.42	-F		-CH ₂ -NMe ₂	17.16	$C_{29}H_{29}FN_4O_2$	485 [M+H] ⁺	216- 220	0.35 (D)
23.43	-F		-CH ₂ -NMe ₂	17.16	$C_{30}H_{31}FN_4O_2$	499 [M+H] ⁺	214- 217	0.35 (D)
23.44	-F		-CH ₂ -NMe ₂	17.16	$C_{31}H_{28}FN_5O_2$	522 [M+H] ⁺	205- 210	0.35 (D)
23.45	-F		-CH ₂ -NMe ₂	17.16	$C_{32}H_{35}FN_4O_2$	527 [M+H] ⁺	235- 237	0.35 (D)
23.46	-F		-CH ₂ -NMe ₂	17.16	$C_{31}H_{28}FN_5O_2$	520 [M-H] ⁻	135- 140	0.20 (D)
23.47	-F		-CH ₂ -NMe ₂	17.16	$C_{29}H_{31}FN_4O_2$	487 [M+H] ⁺	210- 215	0.20 (D)

23.48	-F		-CH ₂ -NMe ₂	17.16	C ₃₀ H ₃₃ FN ₄ O ₂	501 [M+H] ⁺	202- 206	0.25 (D)
23.49	-F		-CH ₂ -NMe ₂	17.16	C ₃₃ H ₃₇ FN ₄ O ₂	541 [M+H] ⁺	198- 203	0.35 (D)
23.50	-F		-CH ₂ -NMe ₂	17.16	C ₂₈ H ₂₉ FN ₄ O ₃	489 [M+H] ⁺	173- 177	0.35 (D)
23.51	-F		-CH ₂ -NMe ₂	17.16	C ₃₃ H ₃₁ FN ₄ O ₃	549 [M-H] ⁻	202- 207	0.50 (C)
23.52	-F		-CH ₂ -NMe ₂	17.16	C ₃₁ H ₃₅ FN ₄ O ₂	513 [M-H] ⁻	203- 209	0.45 (C)
23.53	-F		-CH ₂ -NMe ₂	17.16	C ₃₀ H ₂₇ FN ₄ O ₂ S	527 [M+H] ⁺	245- 250	0.35 (C)
23.54	-F		-CH ₂ -NMe ₂	17.16	C ₃₀ H ₃₃ FN ₄ O ₂	501 [M+H] ⁺	248- 252	0.45 (C)

23.55	-F		-CH ₂ -NMe ₂	17.16	C ₃₀ H ₂₇ FN ₄ O ₃	511 [M+H] ⁺	216- 219	0.30 (C)
23.56	-F		-CH ₂ -NMe ₂	17.16	C ₃₁ H ₂₈ FN ₅ O ₂	522 [M+H] ⁺	167- 170	0.20 (D)

*Eluant mixtures:

(A): silica gel, methylene chloride/ethanol/ammonia = 20:1:0.01

(B): silica gel, methylene chloride/methanol/ammonia = 9:1:0.01

(C): aluminium oxide, methylene chloride/methanol = 19:1

(D): silica gel, methylene chloride/methanol/ammonia = 9:1:0.1

(E): silica gel, methylene chloride/methanol/ammonia = 8:2:0.2

(F): aluminium oxide, methylene chloride/methanol = 9:1

The following were alternatively used as acylating agents:

benzoylchloride, propionylchloride, phenylacetylchloride, cyclopropanecarbonylchloride, cyclobutanecarbonylchloride, pyridin-2-yl-carbonylchloride, pyridin-3-yl-carbonylchloride, pyridin-4-yl-carbonylchloride, cyclohexylcarbonylchloride, isobutyrylchloride, 3-methylbutyrylchloride, cyclohexylmethylcarbonylchloride, methoxyacetylchloride, 2-methoxy-benzoylchloride, tert.-butylacetylchloride, thiophen-2-carbonylchloride, pivaloylchloride, 2-furoyl-chloride.

Example 24.0

3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-phenylsulphonylaminomethyl-phenyl)-methylene]-6-chloro-2-indolinone

100 mg of 3-Z-[1-(4-dimethylaminomethyl-anilino)-1-(4-aminomethyl-phenyl)-methylene]-6-chloro-2-indolinone (educt 14.0) are dissolved in 5 ml methylene chloride and at 0 °C 5 ml of pyridine and 45 µl of benzenesulphonylchloride

are added. The mixture is stirred for 10 minutes at 0 °C and then for 2 hours at ambient temperature. After this time the solvent is eliminated, the residue is suspended in 1N sodium hydroxide solution, suction filtered and washed with a little water. The product is dried at 100°C.

Yield: 87 mg (66% of theory),

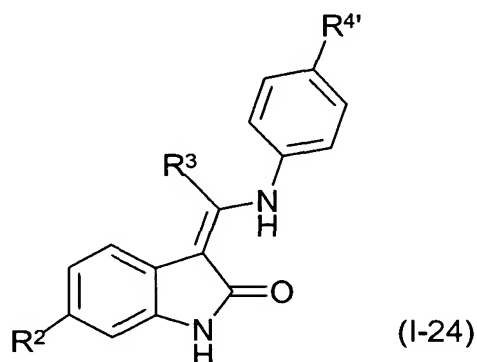
R_f value: 0.30 (silica gel, methylene chloride/methanol/ammonia = 9:1:0.1)

M.p. 170 °C (decomp.)

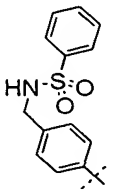
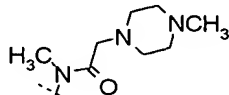
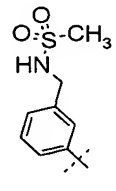
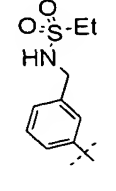
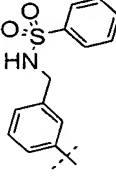
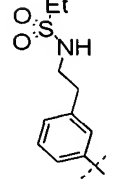
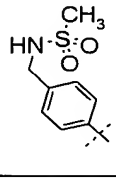
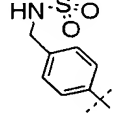
C₃₁H₂₉ClN₄O₃S

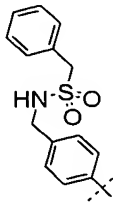
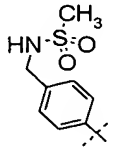
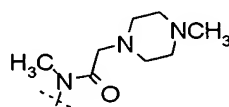
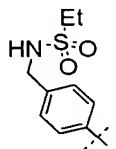
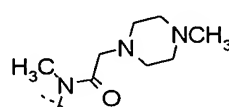
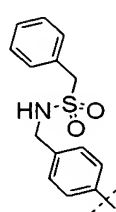
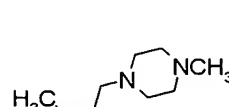
Mass spectrum: m/z = 573/575 [M+H]⁺

The following compounds of general formula I-24 are prepared analogously to Example 24.0:



Example	R ²	R ³	R ^{4'}	educt	empirical formula	mass spectrum	m.p. [°C]	R _f value*
24.1	-Cl			15.0 **	C ₃₁ H ₃₅ ClN ₆ O ₄ S	621/623 [M-H] ⁻	260- 263	0.20 (A)
24.2	-Cl		-CH ₂ -NMe ₂	14.0 **	C ₂₆ H ₂₇ ClN ₄ O ₃ S	511/513 [M+H] ⁺	n. d.	0.35 (A)

24.3	-Cl			15.0	$C_{36}H_{37}ClN_6O_4S$	683/685 [M-H] ⁻	248- 251	0.35 (A)
24.4	-F		-CH ₂ -NMe ₂	17.16 **	$C_{26}H_{27}FN_4O_3S$	495 [M+H] ⁺	170- 180	0.30 (B)
24.5	-F		-CH ₂ -NMe ₂	17.16 ***	$C_{26}H_{27}FN_4O_3S$	509 [M+H] ⁺	200- 204	0.40 (C)
24.6	-F		-CH ₂ -NMe ₂	17.16	$C_{31}H_{29}FN_4O_3S$	557 [M+H] ⁺	125- 130	0.30 (B)
24.7	-F		-CH ₂ -NMe ₂	17.17 ***	$C_{28}H_{31}FN_4O_3S$	521 [M-H] ⁻	100- 110	0.35 (B)
24.8	-F		-CH ₂ -NMe ₂	17.18 **	$C_{26}H_{27}FN_4O_3S$	493 [M-H] ⁻	80- 85	0.35 (B)
24.9	-F		-CH ₂ -NMe ₂	17.18 ***	$C_{27}H_{29}FN_4O_3S$	507 [M-H] ⁻	90- 100	0.40 (B)

24.10	-F		-CH ₂ -NMe ₂	17.18 ****	C ₃₂ H ₃₁ FN ₄ O ₃ S	571 [M+H] ⁺	115- 120	0.35 (B)
24.11	-F			17.19 **	C ₃₁ H ₃₅ FN ₆ O ₄ S	605 [M-H] ⁻	205- 210	0.25 (B)
24.12	-F			17.19 ***	C ₃₂ H ₃₇ FN ₆ O ₄ S	619 [M-H] ⁻	212- 215	0.25 (B)
24.13	-F			17.19 ****	C ₃₇ H ₃₉ FN ₆ O ₄ S	683 [M+H] ⁺	115- 120	0.35 (B)

*Eluant mixtures:

(A): silica gel, methylene chloride/methanol/ammonia = 9:1:0.01

(B): silica gel, methylene chloride/methanol/ammonia = 9:1:0.1

(C): aluminium oxide, methylene chloride/methanol = 19:1

** with methanesulphonylchloride as sulphonating agent

*** with ethanesulphonylchloride as sulphonating agent

**** with α -toluenesulphonylchloride as sulphonating agent

Example 25.0

3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone x ethanesulphonic acid

a) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

10.25 g of 1-acetyl-3-(1-ethoxy-1-phenylmethylene)-6-chloro-2-indolinone (educt IX) and 8.6 g of N-[(4-methyl-piperazin-1-yl)-methylcarbonyl]-N-methyl-p-phenylenediamine (educt XV.204) are dissolved in 100 ml of dimethylformamide and stirred for 4 hours at 120°C. After cooling, 20 ml of 6 N sodium hydroxide solution are added and the mixture is stirred for another hour at ambient temperature. Water is added, the precipitate formed is suction filtered and washed with a little water and 200 ml of ethanol. The residue is dissolved in methylene chloride, extracted with water and dried over sodium sulphate. After the solvent has been eliminated the substance is again washed with a little methanol and dried in vacuo at 100 °C.

Yield: 15.48 g (74% of theory),

R_f value: 0.50 (silica gel, methylene chloride/methanol/ammonia 5:1:0.01)

IR spectrum: 1645 cm⁻¹

M.p. 265-269 °C

C₂₉H₃₀ClN₅O₂

Mass spectrum: m/z = 515/517 [M]⁺

Elemental analysis: calculated:	C 67.49	H 5.86	Cl 6.87	N 13.57
found:	C 67.42	H 5.83	Cl 6.97	N 13.59

b) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone x ethanesulphonic acid

1.5 g of 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone are placed in 22.5 ml of methanol, the mixture is heated to 50°C and 0.25 ml of ethanesulphonic acid in 0.14 g water are added dropwise. The mixture is slowly cooled to

ambient temperature, lastly by means of an ice bath to 0°C. The precipitate formed is suction filtered and washed with a little tert.butylmethylether. The residue is dried in vacuo at 40°C.

Yield: 1.7 g (93% of theory),

IR spectrum: 1655 cm⁻¹

M.p. 307 °C

$C_{29}H_{30}ClN_5O_2 \times C_2H_6O_3S$

Mass spectrum: m/z = 516/518 [M+H]⁺

Example 26

Dry ampoule containing 75 mg of active substance per 10 ml

Composition:

Active substance	75.0 mg
Mannitol	50.0 mg
water for injections	ad 10.0 ml

Preparation:

Active substance and mannitol are dissolved in water. After packaging, the solution is freeze-dried. To produce the solution ready for use, the product is dissolved in water for injections.

Example 27

Dry ampoule containing 35 mg of active substance per 2 ml

Composition:

Active substance	35.0 mg
Mannitol	100.0 mg
water for injections	ad 2.0 ml

Preparation:

Active substance and mannitol are dissolved in water. After packaging, the solution is freeze-dried.

To produce the solution ready for use, the product is dissolved in water for injections.

Example 28

Tablet containing 50 mg of active substance

Composition:

(1) Active substance	50.0 mg
(2) Lactose	98.0 mg
(3) Maize starch	50.0 mg
(4) Polyvinylpyrrolidone	15.0 mg
(5) Magnesium stearate	<u>2.0 mg</u>
	215.0 mg

Preparation:

(1), (2) and (3) are mixed together and granulated with an aqueous solution of (4). (5) is added to the dried granulated material. From this mixture tablets are pressed, biplanar, faceted on both sides and with a dividing notch on one side.

Diameter of the tablets: 9 mm.

Example 29Tablet containing 350 mg of active substance

Composition:

(1) Active substance	350.0 mg
(2) Lactose	136.0 mg
(3) Maize starch	80.0 mg
(4) Polyvinylpyrrolidone	30.0 mg
(5) Magnesium stearate	<u>4.0 mg</u>
	600.0 mg

Preparation

(1), (2) and (3) are mixed together and granulated with an aqueous solution of (4). (5) is added to the dried granulated material. From this mixture tablets are pressed, biplanar, faceted on both sides and with a dividing notch on one side.

Diameter of the tablets: 12 mm.

Example 30Capsules containing 50 mg of active substance

Composition:

(1) Active substance	50.0 mg
(2) Dried maize starch	58.0 mg
(3) Powdered lactose	50.0 mg
(4) Magnesium stearate	<u>2.0 mg</u>
	160.0 mg

Preparation:

(1) is triturated with (3). This trituration is added to the mixture of (2) and (4) with vigorous mixing.

This powder mixture is packed into size 3 hard gelatine capsules in a capsule filling machine.

Example 31

Capsules containing 350 mg of active substance

Composition:

(1) Active substance	350.0 mg
(2) Dried maize starch	46.0 mg
(3) Powdered lactose	30.0 mg
(4) Magnesium stearate	<u>4.0 mg</u>
	430.0 mg

Preparation:

(1) is triturated with (3). This trituration is added to the mixture of (2) and (4) with vigorous mixing.

This powder mixture is packed into size 0 hard gelatine capsules in a capsule filling machine.

Example 32

Suppositories containing 100 mg of active substance

1 suppository contains:

Active substance	100.0 mg
Polyethyleneglycol (M.W. 1500)	600.0 mg
Polyethyleneglycol (M.W. 6000)	460.0 mg
Polyethylenesorbitan monostearate	<u>840.0 mg</u>
	2000.0 mg

Preparation:

The polyethyleneglycol is melted together with polyethylene sorbitan monostearate. At 40°C the ground active substance is homogeneously

dispersed in the melt. It is cooled to 38°C and poured into slightly chilled suppository moulds.

The following compounds may be prepared analogously to the foregoing

Examples:

- (1) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (2) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-ethyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (3) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (4) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (5) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-ethyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (6) 3-Z-[1-(3-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (7) 3-Z-[1-(3-(N-(2-dimethylamino-ethyl)-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolino
- (8) 3-Z-[1-(3-(N-(2-dimethylamino-ethyl)-N-ethyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolino
- (9) 3-Z-[1-(3-(N-(3-dimethylamino-propyl)-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (10) 3-Z-[1-(3-(N-(3-dimethylamino-propyl)-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (11) 3-Z-[1-(3-(N-(3-dimethylamino-propyl)-N-ethyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (12) 3-Z-[1-(3-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (13) 3-Z-[1-(3-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (14) 3-Z-[1-(4-methylsulphonyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (15) 3-Z-[1-anilino-1-phenyl-methylene]-6-chloro-2-indolinone
- (16) 3-Z-[1-(4-nitro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (17) 3-Z-[1-(4-fluoro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (18) 3-Z-[1-(4-chloro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (19) 3-Z-[1-(4-bromo-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (20) 3-Z-[1-(4-iodo-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (21) 3-Z-[1-(4-cyano-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (22) 3-Z-[1-(4-methoxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (23) 3-Z-[1-(4-ethoxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (24) 3-Z-[1-(4-trifluoromethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (25) 3-Z-[1-(4-methylmercapto-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (26) 3-Z-[1-(4-(isopropylaminomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (27) 3-Z-[1-(4-(anilinomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (28) 3-Z-[1-(4-(isobutylaminomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (29) 3-Z-[1-(4-(cyclohexylaminomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (30) 3-Z-[1-(4-(benzylaminomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (31) 3-Z-[1-(4-((N-isopropyl-N-methyl-amino)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (32) 3-Z-[1-(4-((N-ethyl-N-propyl-amino)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (33) 3-Z-[1-(4-((N-ethyl-N-isopropyl-amino)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (34) 3-Z-[1-(4-(dipropylaminomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (35) 3-Z-[1-(4-(diisopropylaminomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (36) 3-Z-[1-(4-((N-benzyl-N-ethyl-amino)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (37) 3-Z-[1-(4-(dibenzylaminomethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (38) 3-Z-[1-(4-(3,6-dihydro-2H-pyridine-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (39) 3-Z-[1-(4-(3,5-dimethyl-piperidin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (40) 3-Z-[1-(4-(azepan-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (41) 3-Z-[1-(4-(2-amino-ethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (42) 3-Z-[1-(4-(2-methylamino-ethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (43) 3-Z-[1-(4-(2-ethylamino-ethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (44) 3-Z-[1-(4-(2-diethylamino-ethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (45) -Z-[1-(4-(2-piperidin-1-yl-ethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (46) 3-Z-[1-(4-(2-acetylamino-ethyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (47) 3-Z-[1-(4-(3-amino-propyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (48) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (49) 3-Z-[1-(4-(N-ethylaminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (50) 3-Z-[1-(4-(N-diethylaminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (51) 3-Z-[1-(4-(N-dipropylaminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (52) 3-Z-[1-(4-(N-((N-ethyl-N-methyl-amino)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (53) 3-Z-[1-(4-(N-((N-ethyl-N-propyl-amino)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (54) 3-Z-[1-(4-(N-((N-methyl-N-propyl-amino)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (55) 3-Z-[1-(4-(N-dimethylaminomethylcarbonyl-N-ethyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (56) 3-Z-[1-(4-(N-dimethylaminomethylcarbonyl-N-propyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (57) 3-Z-[1-(4-(N-dimethylaminomethylcarbonyl-N-butyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (58) 3-Z-[1-(4-(N-(2-amino-ethylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (59) 3-Z-[1-(4-(N-(2-diethylamino-ethylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (60) 3-Z-[1-(4-(N-acetyl-N-(2-aminoethyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (61) -Z-[1-(4-(N-acetyl-N-(2-methylamino-ethyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (62) 3-Z-[1-(4-(N-acetyl-N-(3-amino-propyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (63) 3-Z-[1-(4-(N-acetyl-N-(3-methylamino-propyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (64) 3-Z-[1-(4-(N-acetyl-N-(2-piperidin-1-yl-ethyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (65) 3-Z-[1-(4-(N-acetyl-N-(aminocarbonylmethyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (66) 3-Z-[1-(4-(N-acetyl-N-(piperidin-1-yl-carbonylmethyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (67) 3-Z-[1-(4-(N-(2-ethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (68) 3-Z-[1-(4-(N-(2-diethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (69) 3-Z-[1-(4-(N-(2-pyrrolidin-1-yl-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (70) 3-Z-[1-(4-(N-(2-piperidin-1-yl-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (71) 3-Z-[1-(4-(N-(2-piperazin-1-yl-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (72) 3-Z-[1-(4-(N-(2-(4-morpholin-1-yl)-ethyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (73) 3-Z-[1-(4-(2-diethylamino-ethoxy)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (74) 3-Z-[1-(4-(3-dimethylamino-propoxy)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (75) 3-Z-[1-(4((N-phenethyl-N-methyl-amino)-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (76) 3-Z-[1-(4-carbamoylmethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (77) 3-Z-[1-(4-methylcarbamoylmethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (78) 3-Z-[1-(4-dimethylcarbamoylmethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (79) 3-Z-[1-(4-tetrazol-5-yl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (80) 3-Z-[1-(4-((2-methoxy-ethyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (81) 3-Z-[1-(4-(di-(2-methoxy-ethyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (82) 3-Z-[1-(4-((N-tert.butoxycarbonyl-2-amino-ethyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (83) 3-Z-[1-(4-((N-tert.butoxycarbonyl-3-amino-propyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (84) 3-Z-[1-(4-((2-amino-ethyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (85) 3-Z-[1-(4-((3-amino-propyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (86) 3-Z-[1-(4-((2-acetyl-amino-ethyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (87) 3-Z-[1-(4-((3-acetylamino-propyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (88) 3-Z-[1-(4-((2-methylsulphonylamino-ethyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (89) 3-Z-[1-(4-((3-methylsulphonylamino-propyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (90) 3-Z-[1-(4-(N-(2-amino-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (91) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (92) 3-Z-[1-(4-(N-(2-acetylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (93) 3-Z-[1-(4-(N-(2-methylsulphonylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (94) 3-Z-[1-(4-(carbamoylmethyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (95) 3-Z-[1-(4-(dimethylcarbamoyl-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (96) 3-Z-[1-(4-(methylcarbamoyl-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (97) 3-Z-[1-(4-(N-(pyridin-4-yl-methylcarbonyl)-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (98) 3-Z-[1-(4-(N-(4-propyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (99) 3-Z-[1-(4-(N-(4-butyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (100) 3-Z-[1-(4-(N-(4-isopropyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (101) 3-Z-[1-(4-(N-(4-(2-hydroxy-ethyl)-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (102) 3-Z-[1-(4-(N-(4-(2-methoxy-ethyl)-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (103) 3-Z-[1-(4-(N-(4-(2-ethoxy-ethyl)-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (104) 3-Z-[1-(4-(N-(4-(2-amino-ethyl)-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (105) 3-Z-[1-(4-(N-(4-(2-dimethylamino-ethyl)-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (106) 3-Z-[1-(4-(N-(4-(2-phenyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (107) 3-Z-[1-(4-(N-(4-(5-methyl-2,5-diaza-bicyclo[2.2.1]-hept-2-yl)-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (108) 3-Z-[1-(4-(N-(homopiperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (109) 3-Z-[1-(4-(N-(3,4,5-trimethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (110) 3-Z-[1-(4-(N-(2,4,6-trimethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (111) 3-Z-[1-(4-(N-(trans-2,4,5-trimethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (112) 3-Z-[1-(4-(N-(cis-2,4,5-trimethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (113) 3-Z-[1-(4-(N-(2,4-dimethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (114) 3-Z-[1-(4-(N-(3,4-dimethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (115) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-ethyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (116) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-propyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (117) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-isopropyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (118) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-butyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (119) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-phenyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (120) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-benzyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (121) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-aminocarbonylmethyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (122) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-dimethylaminocarbonylmethyl-amino)-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (123) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-bromo-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (124) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-chloro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (125) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-fluoro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (126) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-methyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (127) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-methoxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (128) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-hydroxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (129) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-cyano-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (130) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-trifluormethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (131) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-aminocarbonyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (132) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-amino-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (133) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-nitro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (134) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-acetylamino-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (135) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3-methylsulphonylamino-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (136) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3,5-dibromo-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (137) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-3,5-dichloro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (138) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-5-nitro-6-chloro-2-indolinone
- (139) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-(pyridin-3-yl-amino))-1-phenyl-methylene]-6-chloro-2-indolinone
- (140) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-(pyridin-2-yl-amino))-1-phenyl-methylene]-6-chloro-2-indolinone
- (141) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-bromo-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (142) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-chloro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (143) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-fluoro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (144) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-methyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (145) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-methoxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (146) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-hydroxy-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (147) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-cyano-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (148) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-trifluormethyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (149) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-aminocarbonyl-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (150) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-amino-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone

- (151) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-nitro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (152) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-acetyl-amino-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (153) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3-methylsulphonylamino-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (154) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3,5-dibromo-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (155) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-3,5-dichloro-anilino)-1-phenyl-methylene]-6-chloro-2-indolinone
- (156) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-chloro-2-indolinone
- (157) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-chloro-2-indolinone
- (158) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-chloro-2-indolinone
- (159) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-chloro-2-indolinone
- (160) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-chloro-2-indolinone
- (161) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-chloro-2-indolinone
- (162) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-chloro-2-indolinone
- (163) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone
- (164) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone
- (165) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone

- (166) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone
- (167) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone
- (168) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone
- (169) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone
- (170) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone
- (171) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-chloro-2-indolinone
- (172) 3-Z-[1-(4-(N-(pyridin-4-yl-methylcarbonyl)-amino)-anilino)-1-phenyl-methylene]-6-cyano-2-indolinone
- (173) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-cyano-2-indolinone
- (174) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-cyano-2-indolinone
- (175) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-cyano-2-indolinone
- (176) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-cyano-2-indolinone
- (177) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-cyano-2-indolinone
- (178) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-cyano-2-indolinone
- (179) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-cyano-2-indolinone
- (180) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-cyano-2-indolinone

- (181) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3,4-dimethoxy-phenyl)-methylene]-6-cyano-2-indolinone
- (182) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-ethylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (183) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propionyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (184) 3-Z-[1-(4-(N-(3-dimethylaminopropyl)-N-methylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (185) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-butyryl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (186) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-isobutyryl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (187) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (188) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-benzoyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (189) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-phenylacetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (190) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(pyrid-3-yl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (191) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(furan-2-yl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (192) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(2-methoxy-benzoyl)-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (193) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-isopropylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (194) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-benzylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (195) 3-Z-[1-(4-(N-(4-benzyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (196) 3-Z-[1-(4-(N-(piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (197) 3-Z-[1-(4-(N-(morpholin-4-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone

- (198) 3-Z-[1-(4-(N-(piperidin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (199) 3-Z-[1-(4-(N-(benzylmethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (200) 3-Z-[1-(4-(N-(methylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (201) 3-Z-[1-(4-((2,6-dimethyl-piperidin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (202) 3-Z-[1-(4-((N-(2-(2-methoxy-ethoxy)-ethyl)-N-methyl-amino)-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (203) 3-Z-[1-(4-(triazol-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (204) 3-Z-[1-(4-(di-(2-hydroxy-ethyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (205) 3-Z-[1-(4-(5-methyl-imidazol-4-yl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (206) 3-Z-[1-(4-(morpholin-4-yl-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (207) 3-Z-[1-(4-(N-benzyl-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (208) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-butylsulphonyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (209) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methoxyacetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (210) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(3,4-dimethoxy-benzoyl)-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (211) 3-Z-[1-(4-(N-(2-hydroxy-ethyl)-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (212) 3-Z-[1-(4-(N-(2-benzylmethylamino-ethyl)-N-propionyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (213) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-propionyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (214) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-(pyrid-4-yl-carbonyl)-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone

- (215) 3-Z-[1-(4-(N-(phthalimido-2-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (216) 3-Z-[1-(4-(N-aminomethylcarbonyl-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (217) 3-Z-[1-(4-(N-tert.butoxycarbonyl-methylamino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (218) 3-Z-[1-(4-(N-(di-(2-hydroxy-ethyl)-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (219) 3-Z-[1-(4-(N-(pyrrolidin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (220) 3-Z-[1-(4-(N-(imidazol-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (221) 3-Z-[1-(4-(N-(2-(4-tert.butoxycarbonyl-piperazin-1-yl)-ethylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (222) 3-Z-[1-(4-(N-(2-(piperazin-1-yl)-ethylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (223) 3-Z-[1-(4-(N-(2-benzylmethylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (224) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (225) 3-Z-[1-(4-((4-tert.butoxycarbonyl-piperazin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (226) 3-Z-[1-(4-(piperazin-1-yl-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (227) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-3-methoxy-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (228) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (229) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-pyridin-2-yl-amino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (230) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-(pyridin-3-yl-amino))-1-phenyl-methylene]-6-fluoro-2-indolinone
- (231) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-3-methyl-pyrrol-3-yl-amino)-1-phenyl-methylene]-6-fluoro-2-indolinone

- (232) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl)-3-methyl-pyrrol-3-yl-amino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (233) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (234) 3-Z-[1-(4-(1-(2-dimethylamino-ethyl)-imidazol-2-yl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (235) 3-Z-[1-(4-(piperidin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (236) 3-Z-[1-(4-(4-tert.butoxycarbonyl-piperazin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (237) 3-Z-[1-(4-(piperazin-1-yl-carbonyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (238) 3-Z-[1-(4-(N-cyclohexyl-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (239) 3-Z-[1-(4-(N-(2-(4-methyl-piperazin-1-yl)-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (240) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (241) 3-Z-[1-(4-((4-dimethylamino-piperidin-1-yl)-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (242) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (243) 3-Z-[1-(4-(N-tert.butoxycarbonyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (244) 3-Z-[1-(4-aminomethyl-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (245) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-isopropyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (246) 3-Z-[1-(4-(N-(4-tert.butoxycarbonyl-piperazin-1-yl-methylcarbonyl)-N-isopropyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (247) 3-Z-[1-(4-(N-(piperazin-1-yl-methylcarbonyl)-N-isopropyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (248) 3-Z-[1-(4-(diethylamino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone

- (249) 3-Z-[1-(4-(N-propyl-N-methyl-amino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (250) 3-Z-[1-(4-(4-methyl-piperazin-1-yl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (251) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (252) 3-Z-[1-(4-(N-(4-tert.butoxycarbonyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (253) 3-Z-[1-(4-(N-(piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (254) 3-Z-[1-(4-(N-(hydroxy-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (255) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (256) 3-Z-[1-(4-(N-(N-(tert.butoxycarbonyl-3-amino-propyl)-N-methyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (257) 3-Z-[1-(4-(N-(N-(3-amino-propyl)-N-methyl)-amino-methyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (258) 3-Z-[1-(4-(N-(4-methyl-homopiperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (259) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-3-cyano-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (260) 3-Z-[1-(4-(N-(4-ethyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (261) 3-Z-[1-(4-(N-(1-methyl-piperidin-4-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (262) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (263) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methyl-carbamoylmethyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (264) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-carbamoylmethyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (265) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonylmethyl)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone

- (266) 3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (267) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-aminocarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (268) 3-Z-[1-(4-(N-(1-methyl-piperidin-4-yl-aminocarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (269) 3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (270) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (271) 3-Z-[1-(4-(N-(N-(3-dimethylamino-propyl)-aminocarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (272) 3-Z-[1-(4-(N-(pyridin-4-yl-methylaminocarbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (273) 3-Z-[1-(4-(N-(1-methyl-piperidin-4-oxy-carbonyl)-N-methyl-amino)-anilino)-1-phenyl-methylene]-6-fluoro-2-indolinone
- (274) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (275) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (276) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (277) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (278) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (279) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (280) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-oxo-oxazolidin-3-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (281) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone

- (282) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone
- (283) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone
- (284) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone
- (285) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone
- (286) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone
- (287) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone
- (288) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-amino)-N-methyl-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone
- (289) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-methoxycarbonylamino-phenyl)-methylene]-6-fluoro-2-indolinone
- (290) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (291) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (292) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (293) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (294) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (295) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (296) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-fluoro-2-indolinone

- (297) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-(2-oxo-pyrrolidine-2-yl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (298) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (299) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (300) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (301) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (302) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (303) 3-Z-[1-(4-(N-(3-methylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (304) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (305) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (306) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (307) 3-Z-[1-(4-methylaminomethyl-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (308) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (309) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (310) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (311) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (312) 3-Z-[1-(4-aminomethyl-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (313) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone

- (314) 3-Z-[1-(3-(methylamino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (315) 3-Z-[1-(3-(2-dimethylamino-ethyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (316) 3-Z-[1-(3-(3-dimethylamino-propyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (317) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (318) 3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (319) 3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (320) 3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (321) 3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (322) 3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (323) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (324) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (325) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (326) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (327) 3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (328) 3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (329) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone

- (330) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (331) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (332) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (333) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (334) 3-Z-[1-(4-(N-(3-methylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (335) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (336) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (337) 3-Z-[1-(4-methylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (338) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (339) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (340) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (341) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (342) 3-Z-[1-(4-aminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (343) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (344) 3-Z-[1-(3-(methylamino-methyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (345) 3-Z-[1-(3-(2-dimethylamino-ethyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (346) 3-Z-[1-(3-(3-dimethylamino-propyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone

- (347) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (348) 3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (349) 3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (350) 3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (351) 3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (352) 3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (353) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (354) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (355) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (356) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (357) 3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (358) 3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-chloro-2-indolinone
- (359) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (360) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (361) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (362) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone

- (363) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (364) 3-Z-[1-(4-(N-(3-methylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (365) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (366) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (367) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (368) 3-Z-[1-(4-methylaminomethyl-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (369) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (370) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (371) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (372) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (373) 3-Z-[1-(4-aminomethyl-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (374) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (375) 3-Z-[1-(3-(methylamino-methyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (376) 3-Z-[1-(3-(2-dimethylamino-ethyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (377) 3-Z-[1-(3-(3-dimethylamino-propyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (378) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (379) 3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone

- (380) 3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (381) 3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (382) 3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (383) 3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (384) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (385) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (386) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (387) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (388) 3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (389) 3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (390) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (391) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (392) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (393) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (394) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (395) 3-Z-[1-(4-(N-(3-methylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone

- (396) 3-Z-[1-(4-(dimethylamino-methyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (397) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (398) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (399) 3-Z-[1-(4-methylaminomethyl-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (400) 3-Z-[1-(4-(N-(4-methyl-piperazin-1-yl-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (401) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (402) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (403) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (404) 3-Z-[1-(4-aminomethyl-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (405) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (406) 3-Z-[1-(3-(methylamino-methyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (407) 3-Z-[1-(3-(2-dimethylamino-ethyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (408) 3-Z-[1-(3-(3-dimethylamino-propyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (409) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (410) 3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (411) 3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (412) 3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone

- (413) 3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (414) 3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (415) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (416) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (417) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (418) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (419) 3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (420) 3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-chloro-2-indolinone
- (421) 3-Z-[1-(4-(N-(dimethylamino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (422) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (423) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (424) 3-Z-[1-(4-(N-(3-methylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (425) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (426) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (427) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (428) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

- (429) 3-Z-[1-(4-aminomethyl-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (430) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (431) 3-Z-[1-(3-(methylamino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (432) 3-Z-[1-(3-(2-dimethylamino-ethyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (433) 3-Z-[1-(3-(3-dimethylamino-propyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (434) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (435) 3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (436) 3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (437) 3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (438) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (439) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (440) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (441) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (442) 3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (443) 3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (444) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

- (445) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (446) 3-Z-[1-(4-(N-(3-methylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (447) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (448) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (449) 3-Z-[1-(4-methylaminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (450) 3-Z-[1-(4-(4-methyl-piperazin-1-yl-carbonyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (451) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (452) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (453) 3-Z-[1-(4-aminomethyl-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (454) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (455) 3-Z-[1-(3-(methylamino-methyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (456) 3-Z-[1-(3-(2-dimethylamino-ethyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (457) 3-Z-[1-(3-(3-dimethylamino-propyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (458) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (459) 3-Z-[1-(4-(N-(dimethylamino-carbonylmethyl)-N-methylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (460) 3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone

- (461) 3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (462) 3-Z-[1-(4-(1-methyl-imidazol-2-yl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (463) 3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (464) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (465) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (466) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (467) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (468) 3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (469) 3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-(2-carboxy-ethyl)-phenyl)-methylene]-6-fluoro-2-indolinone
- (470) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (471) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (472) 3-Z-[1-(4-(N-(3-methylamino-propyl)-N-acetyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (473) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (474) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (475) 3-Z-[1-(4-methylaminomethyl-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (476) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone

- (477) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (478) 3-Z-[1-(4-aminomethyl-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (479) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (480) 3-Z-[1-(3-(methylamino-methyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (481) 3-Z-[1-(3-(2-dimethylamino-ethyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (482) 3-Z-[1-(3-(3-dimethylamino-propyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (483) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (484) 3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (485) 3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (486) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (487) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (488) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (489) 3-Z-[1-(4-(N-(4-dimethylamino-butyl-carbonyl)-N-methyl-amino)-anilino)-1-(4-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (490) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (491) 3-Z-[1-(4-(N-(2-methylamino-ethyl)-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (492) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone

- (493) 3-Z-[1-(4-(N-(3-methylamino-propyl)-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (494) 3-Z-[1-(4-(3-dimethylamino-propyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (495) 3-Z-[1-(4-ethylaminomethyl-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (496) 3-Z-[1-(4-methylaminomethyl-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (497) 3-Z-[1-(4-(N-(3-dimethylamino-propyl)-N-methylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (498) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-propylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (499) 3-Z-[1-(4-aminomethyl-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (500) 3-Z-[1-(3-(dimethylamino-methyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (501) 3-Z-[1-(3-(methylamino-methyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (502) 3-Z-[1-(3-(2-dimethylamino-ethyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (503) 3-Z-[1-(3-(3-dimethylamino-propyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (504) 3-Z-[1-(4-(2-dimethylamino-ethyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (505) 3-Z-[1-(4-(N-methyl-N-methylsulphonyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (506) 3-Z-[1-(4-(N-methyl-N-acetyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (507) 3-Z-[1-(4-(N-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methylcarbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (508) 3-Z-[1-(4-(2-diethylamino-ethyl-sulphonyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone

- (509) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (510) 3-Z-[1-(4-(N-(2-dimethylamino-ethyl)-N-methyl-amino-methyl)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (511) 3-Z-[1-(4-(2-dimethylamino-ethoxy)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone
- (512) 3-Z-[1-(4-(N-(3-dimethylamino-propyl-carbonyl)-N-methyl-amino)-anilino)-1-(3-carboxymethyl-phenyl)-methylene]-6-fluoro-2-indolinone

In the above Tables

Me denotes methyl,
Et denotes ethyl,
Pr denotes propyl,
nPr denotes n-propyl,
iPr denotes isopropyl,
nBu denotes n-butyl,
tBu denotes tert.-butyl and
Bn denotes benzyl.